

# Setting Priorities for Hazardous Waste Minimization

**Appendices** 



#### TABLE OF CONTENTS

Appendix 1.	Sample BRS Data Politis
Appendix 2.	BRS Code Descriptions
Appendix 3.	Matching Process - WR/GM
Appendix 4.	Documents Used as Sources of Concentration Data
Appendix 5.	Retrieval from GENSUR
Appendix 6.	Waste Characterizations for Top 150 Routinely Generated Combusted Waste Stream Combinations
Appendix 7.	Summary of Prioritization Systems
Appendix 8.	HRS Hazard Data and Pathway Scores from Superfund Chemical Data Matrix
Appendix 9.	States and Regions in Which Top 100 Ranked Wastestream Combinations are Generated

#### APPENDIX 1 SAMPLE BRS DATA FORMS

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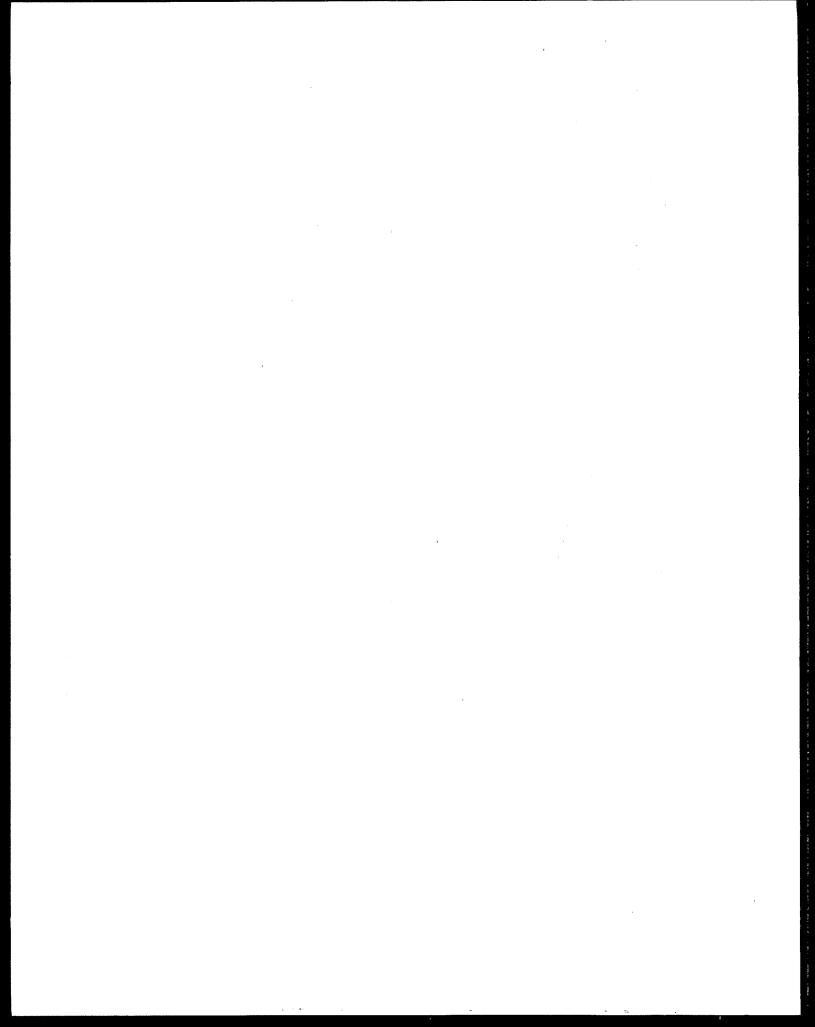
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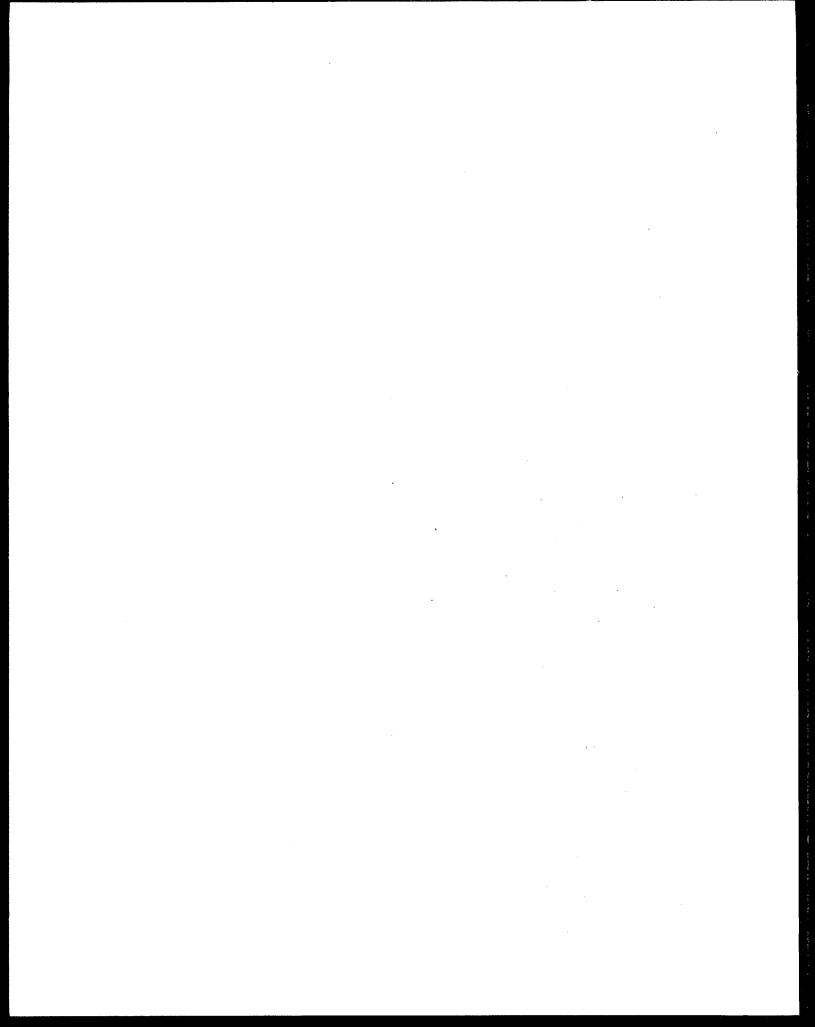
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#### APPENDIX 2 BRS CODE DESCRIPTIONS

EPA Hazardous Waste Codes
Source Codes
Form Codes
SIC Codes
Origin Codes



#### **EPA HAZARDOUS WASTE CODES**

Code	Waste description	Code	Waste description
CHAR	ACTERISTIC HAZARDOUS WASTE	D026	Cresol
D001	Ignitable waste	D027	1,4-Dichlorobenzene
D002	Corrosive waste	D028	1,2-Dichloroethane
D003	Reactive waste	D029 1.1-Dichloroethylens	1,1-Dichloroethylene
D004	Arsenic	.D030	2,4-Dinitrotoluene
D005	Barium	D031	Heptachlor (and its epoxide)
D006	Cadmium	D032	Hexachlorobenzene
D007	Chromium	D033	Hexachlorobutadiene
D007	Lead	D034	Hexachloroethane
D009		D035	Methyl ethyl ketone
	Mercury	D036	Nitrobenzene
D010	Selenium	D037	Pentachlorophenol
D011	Silver	D038	Pyridine
D012	Endrin(1,2,3,4,10,10-hexachloro-1,7-epoxy-1,4,4a,5,6,7,8,8a-octahydro-1,4-endo, endo-	D039	Tetrachloroethylene
	5,8-dimeth-ano-naphthalene)	D040	Trichloroethylene
D013	Lindane (1,2,3,4,5,6-hexa-	D041	2,4,5-Trichlorophenol
	chlorocyclohexane, gamma isomer)	e, gamma isomer) D042 2,4,6-Trichlorophe	2,4,6-Trichlorophenol
D014	Methoxychlor (1,1,1-trichloro-2,2-bis [p-methoxyphenyl] ethane)	D043	Vinyl chloride
D015	Toxaphene (C <sub>10</sub> H <sub>10</sub> Cl <sub>N</sub> , Technical chlorinated camphene, 67-69 percent chlorine)	HAZA SOUR	RDOUS WASTE FROM NONSPECIFIC
D016	2,4-D (2,4-Dichlorophenoxyacetic acid)		
D017	24.5-TP Silvex (24.5- Trichlorophenoxypropionic acid)	F001	The following spent halogenated solvents used in degreasing: Tetrachloroethylene,
D018	Benzene		trichlorethylene, methylene chloride, 1,1,1- trichloroethane, carbon tetrachloride and
D019	Carbon tetrachloride		chlorinated fluorocarbons; all spent
D020	Chlordane		solvent mixtures/blends used in degreasing containing, before use, a total of ten
D021	Chlorobenzene		percent or more (by volume) of one or
D022	Chloroform		more of the above halogenated solvents or
D023	o-Cresol		those solvents listed in F002, F004, and F005; and still bottoms from the recovery
D024	m-Cresol		of these spent solvents and spent solvent
D025	p-Cresol		mixtures.

Code	Waste description	Code	Waste description
F002	The following spent halogenated solvents: Tetrachloroethylene, methylene chloride, trichloroethylene, 1,1,1-trichloroethane, chlorobenzene, 1,1,2-trichloro-1,2,2-trifluoroethane, ortho-dichlorobenzene, trichlorofluoromethane, and 1,1,2, trichloroethane; all spent solvent mixtures/blends containing, before use, a total of ten percent or more (by volume) of one or more of the above halogenated solvents or those solvents listed in F001, F004, and F005; and still bottoms from the	F005	The following spent non-halogenated solvents: toluene, methyl ethyl ketone, carbon disulfide, isobutanol, pyridine, benzene, 2-ethoxyethanol, and 2-nitropropane; all spent solvent mixtures/blends containing, before use, a total of ten percent or more (by volume) of one or more of the above non-halogenated solvents or those solvents listed in F001, F002, or F004; and still bottoms from the recovery of these spent solvents and spent solvent mixtures.
F003	recovery of these spent solvents and spent solvent mixtures.  The following spent non-halogenated solvents: Xylene, acetone, ethyl acetate, ethyl benzene, ethyl ether, methyl isobutyl ketone, n-butyl alcohol, cyclohexanone, and methanol; all spent solvent mixtures/blends containing, before use, only the above spent non-halogenated solvents; and all spent solvent mixtures/blends containing, before use, one or more of the above non-halogenated solvents, and a total of ten percent or more (by volume) of one or more of those solvents listed in F001, F002, F004, and F005; and still bottoms from the recovery of these spent solvents and spent solvent mixtures.  The following spent non-halogenated solvents: cresols, cresylic acid, and nitrobenzene; all spent solvent mixtures/blends containing, before use, a total of ten percent or more (by volume) of one or more of the above non-halogenated solvents or those solvents	F006	Wastewater treatment sludges from electroplating operations except from the following processes: (1) sulfuric acid anodizing of aluminum; (2) tin plating on carbon steel; (3) zinc plating (segregated basis) on carbon steel; (4) aluminum or zinc-aluminum plating on carbon steel; (5) cleaning/stripping associated with tin, zinc, and aluminum plating on carbon steel; and (6) chemical etching and milling of aluminum.
		F007	Spent cyanide plating bath solutions from electroplating operations.
		F008	Plating bath residues from the bottom of plating baths from electroplating operations in which cyanides are used in the process.
F004		F009	Spent stripping and cleaning bath solutions from electroplating operations in which cyanides are used in the process.
		F010	Quenching bath residues from oil baths from metal heat treating operations in which cyanides are used in the process.
	listed in P001, F002, and F005; and still oottoms from the recovery of these spent solvents and spent solvent mixtures.	F011	Spent cyanide solutions from salt bath pot cleaning from metal heat treating operations.
		F012	Quenching wastewater treatment sludges from metal heat treating operations in which cyanides are used in the process.

Code	Waste description	Code	Waste description	
F019	Wastewater treatment sludges from the chemical conversion coating of aluminum except from zirconium phosphating in aluminum can washing when such phosphating is an exclusive conversion coating process.	F024	Process wastes including, but not limited to, distillation residues, heavy ends, tars, and reactor clean-out wastes, from the production of certain chlorinated aliphatic hydrocarbons by free radical catalyzed processes. These chlorinated aliphatic	
F020	Wastes (except wastewater and spent character	hydrocarbons are those having carbon chain lengths ranging from one to and including five, with varying amounts and positions of chlorine substitution. (This listing does not include wastewaters, wastewater treatment sludges, spent catalysts, and wastes listed in Sections 261.31. or 261.32)		
		Condensed light ends, spent filters and filter aids, and spent desiccant wastes from the production of certain chlorinated		
F021	Wastes (except wastewater and spent carbon from hydrogen chloride purification) from the production or manufacturing use (as a reactant, chemical intermediate, or component in a formulating process) of		aliphatic hydrocarbons, by free radical catalyzed processes. These chlorinated aliphatic hydrocarbons are those having carbon chain lengths ranging from one, to and including five, with varying amounts and positions of chlorine substitution.	
	pentachlorophenol, or of intermediates used to produce derivatives.	F026	Wastes (except wastewater and spent carbon from hydrogen chloride	
F022	Wastes (except wastewater and spent carbon from hydrogen chloride purification) from the manufacturing use (as a reactant, chemical intermediate, or component in a formulating process) of tetra-, penta-, or hexachlorobenzenes under alkaline conditions.		purification) from the production of materials on equipment previously used for the manufacturing use (as a reactant, chemical intermediate, or component in a formulating process) of tetra-, penta-, or hexachlorobenzene under alkaline conditions.	
F023	Wastes (except wastewater and spent carbon from hydrogen chloride purification) from the production of materials on equipment previously used for the production or manufacturing use (as a reactant, chemical intermediate, or component in a formulating process) of tri- and tetrachlorophenols. (This listing	F027	Discarded unused formulations containing tri-, tetra-, or pentachlorophenol or discarded unused formulations containing compounds derived from these chlorophenols. (This listing does not include formulations containing hexachlorophene synthesized from prepurified 2,4,5-trichlorophenol as the sole component.)	
does not include wastes from equipment used only for the production or use of hexachlorophene from highly purified 2,4,5-trichlorophenol.)	F028	Residues resulting from the incineration or thermal treatment of soil contaminated with EPA hazardous waste nos. F020, F021, F022, F023, F026, and F027.		

#### **EPA HAZARDOUS WASTE CODES**

generated in oil/water/solids separators;

(Continued)

Code Waste description Code Waste description F032 Wastewaters, process residuals, tanks and impoundments; ditches and preservative drippage, and spent other conveyances; sumps; and storm formulations from wood preserving water units receiving dry weather flow. processes generated at plants that Sludges generated in storm water units currently use, or have previously used, that do not receive dry weather flow. chlorophenolic formulations [except sludges generated in aggressive biological potentially cross-contaminated wastes that treatment units as defined in Section have had the F032 waste code deleted in 261.31(b)(2)(including sludges generated accordance with Section 261.35 (i.e., the in one or more additional units after newly promulgated equipment cleaning or wastewaters have been treated in replacement standards), and where the aggressive biological treatment units). generator does not not resume or initiate and K051 wastes are exempted from this use of chlorophenolic formulations). This listing. listing does not include K001 bortom F038 Petroleum refinery secondary (emulsified) sediment sludge from the treatment of oil/water/solids separation sludge - Any wastewater from wood preserving sludge and/or float generated from the processes that use creosote and/or physical and/or chemical separation of pentachlorophenoL oil/water/solids in process wastewaters F034 Wastewaters, process residuals, and oily cooling wastewaters from preservative drippage, and spent petroleum refineries. Such wastes include, formulations from wood preserving but are not limited to, all sludges and processes generated at plants that use floats generated in induced air flotation creosote formulations. This listing does (IAF) units, tanks and impoundments, and not include K001 bottom sediment sludge all sludges generated in DAF units. from the treatment of wastewater from Sludges generated in stormwater units that wood preserving processes that use do not receive dry weather flow, sludges crecsote and/or pentachlorophenol. generated in aggressive biological treatment units as defined in Section F035 Wastewaters, process residuals, 261.31(b)(2) (including sludges generated preservative drippage, and spent in one or more additional units after formulations from wood preserving wastewaters have been treated in processes generated at plants that use aggressive biological treatment units), inorganic preservatives containing arsenic and F037, K048, and K051 wastes are or chromium. This listing does not include exempted from this listing. K001 bottom sediment sludge from the treatment of wastewater from wood F039 Leachate resulting from the treatment, preserving processes that use creosote storage, or disposal of wastes classified by and/or peatachlorophenol. more than one waste code under Subpart D, or from a mixture of wastes classified F037 Petroleum refinery primary under Subparts C and D of this part. oil/water/solids separation sludge - Any (Leachate resulting from the management sludge generated from the gravitational of one or more of the following EPA separation of oil/water/solids during the Hazardous Wastes and no other hazardous storage or treatment of process wastes retains its hazardous waste code(s): wastewaters and oily cooling wastewaters F020, F021, F022, F023, F026, F027, from petroleum refineries. Such sludges and/or F028.) include, but are not limited to, those

Code	Waste description	Code	Waste description
HAZA SOUR	RDOUS WASTE FROM SPECIFIC CES	K018	Heavy ends from the fractionation column in ethyl chloride production.
K001	Bottom sediment sludge from the treatment of wastewaters from wood	K019	Heavy ends from the distillation of ethylene dichloride production.
	preserving processes that use creosote and/or pentachlorophenol.	K020	Heavy ends from the distillation of vinyl chloride in vinyl chloride monomer
K002	Wastewater treatment sludge from the production of chrome yellow and orange pigments.	K021	production.  Aqueous spent antimony catalyst waste from fluoromethanes production.
K003	Wastewater treatment sludge from the production of molybdate orange pigments.	K022	Distillation bottom tars from the production of phenol/acetone from
K004	Wastewater treatment sludge from the production of zinc yellow pigments.	K023	cumene.  Distillation light ends from the production of phthalic anhydride from naphthalene.
K005	Wastewater treatment sludge from the production of chrome green pigments.	K024	Distillation bottoms from the production-
K006	Wastewater treatment sludge from the production of chrome oxide green pigments (anhydrous and hydrated).	K025	of phthalic anhydride from naphthalene.  Distillation bottoms from the production of nitrobenzene by the nitration of
K007	Wastewater treatment sludge from the production of iron blue pigments.	K026	benzene.  Stripping still tails from the production of
K008	Oven residue from the production of	row.	methyl ethyl pyridines.
٧,000	chrome oxide green pigments.	K027	Centrifuge and distillation residues from toluene diisocyanate production.
K009	Distillation bottoms from the production of acetaldehyde from ethylene.	K028	Spent catalyst from the hydrochlorinator reactor in the production of 1,1,1-
K010	Distillation side cuts from the production of acetaldehyde from ethylene.		trichloroethane.
K011	Bottom stream from the wastewater stripper in the production of acrylonitrile.	K029	Waste from the product steam stripper in the production of 1,1,1-trichloroethane.
K013	Bottom stream from the acetonitrile column in the production of acrylonitrile.	K030	Column bottoms or heavy ends from the combined production of trichloroethylene and perchloroethylene.
K014	Bottoms from the acetonitrile purification column in the production of acrylonitrile.	K031	By-product salts generated in the
K015	Still bottoms from the distillation of benzyl chloride.	K032	production of MSMA and cacodylic acid.  Wastewater treatment sludge from the production of chlordane.
K016	Heavy ends or distillation residues from the production of carbon tetrachloride.	K033	Wastewater and scrub water from the chlorination of cyclopentadiene in the
K017			production of chlordane.

Code	Waste description	Code	Waste description
K034	Filter solids from the filtration of hexachlorocyclopentadiene in the production of chlordane.	K052	Tank bottoms (leaded) from the petroleum refining industry.
K035	Wastewater treatment sludges generated in the production of creosote.	K060	Ammonia still lime sludge from coking operations.
K036	Still bottoms from toluene reclamation distillation in the production of disulfoton.	K061	Emission control dust/sludge from the primary production of steel in electric furnaces.
K037	Wastewater treatment sludges from the production of disulfoton.	K062	Spent pickle liquor generated by steel finishing operations of facilities within the
K038	Wastewater from the washing and stripping of phorate production.		iron and steel industry (SIC Codes 331 and 332).
K039	Filter cake from the filtration of diethylphosphorodithioic acid in the production of phorate.	K064	Acid plant blowdown slurry/sludge resulting from the thickening of blowdown slurry from primary copper production.
K040	Wastewater treatment sludge from the production of phorate.	K065	Surface impoundment solids contained in and dredged from surface impoundments
K041	Wastewater treatment sludge from the production of toxaphene.	K066	at primary lead smelting facilities.  Sludge from treatment of process
K042	Heavy ends or distillation residues from the distillation of tetrachlorobenzene in	****	wastewater and/or acid plant blowdown from primary ziac production.
K043	the production of 2,4,5-T.	K069	Emission control dust/sludge from secondary lead smelting.
2043	2.6-dichlorophenol waste from the production of 2,4-D.	K071	Brine purification muds from the mercury
K044	Wastewater treatment sludges from the manufacturing and processing of explosives.		cell process in chlorine production, in which separately prepurified brine is not used.
K045	Spent carbon from the treatment of wastewater containing explosives.	K073	Chlorinated hydrocarbon waste from the purification step of the diaphragm cell process using graphite anodes in chlorine
K046	Wastewater treatment sludges from the manufacturing, formulation, and loading of lead-based initiating compounds.	K083	production.  Distillation bottoms from aniline
K047	Pink/red water from TNT operations.	K084	production.
K048	Dissolved air flotation (DAF) float from the petroleum refining industry.	NUO-	Wastewater treatment siudges generated during the production of veterinary pharmaceuticals from arsenic or organo-
K049	Slop oil emulsion solids from the petroleum refining industry.	K085	arsenic compounds.  Distillation or fractionation column
K050	Heat exchanger bundle cleaning sludge from the petroleum refining industry.	,	bottoms from the production of chlorobenzenes.
K051	API separator sludge from the petroleum refining industry.		

Code	Waste description	Code	Waste description
K086	Solvent washes and sludges, caustic washes and sludges, or water washes and sludges	K103	Process residues from aniline extraction from the production of aniline.
	from cleaning tubs and equipment used in the formulation of ink from pigments, driers, soaps, and stabilizers containing	K104	Combined wastewater streams generated from nitrobenzene/aniline production.
K087	chromium and lead.  Decanter tank tar sludge from coking	K105	Separated aqueous stream from the reactor product washing step in the production of chlorobenzenes.
K088	operations.  Spent potliners from primary aluminum reduction.	K106	Wastewater treatment sludge from the mercury cell process in chlorine production.
K090	Emission control dust or sludge from ferrochromiumsilicon production.	K107	Column bottoms from product separation from the production of 1,1-
K091	Emission control dust or sludge from ferrochromium production.		dimethylhydrazine (UDMH) from carboxylic acid hydrazines.
K093	Distillation light ends from the production of phthalic anhydride from ortho-xylene.	K108	Condensed column overheads from product separation and condensed reactor
K094	Distillation bottoms from the production of phthalic anhydride from ortho-xylene.		vent gases from the production of 1.1- dimethylhydrazine (UDMH) from carboxylic acid hydrazides.
K095	Distillation bottoms from the production of 1,1,1-trichloroethane.	K109	Spent filter cartridges from product purification from the production of 1,1-
K096	Heavy ends from the heavy ends column from the production of 1,1,1-trichloroethane.		dimethylhydrazine (UDMH) from carboxylic acid hydrazides.
K097	Vacuum stripper discharge from the chlordane chlorinator in the production of chlordane.	K110	Condensed column overheads from intermediate separation from the production of 1,1-dimethylhydrazine (UDMH) from carboxylic acid hydrazides.
K098	Untreated process wastewater from the production of toxaphene.	K111	Product washwaters from the production of dinitrotoluene via nitration of toluene.
K099	Untreated wastewater from the production of 2,4-D.	K112	Reaction by-product water from the drying column in the production of
K100	Waste leaching solution from acid leaching of emission control dust/sludge from		toluenediamine via hydrogenation of dinitrotoluene.
K101	secondary lead smelting.  Distillation tar residues from the distillation of aniline-based compounds in the production of veterinary	K113	Condensed liquid light ends from purification of toluenediamine in production of toluenediamine via hydrogenation of dinitrotoluene.
	pharmaceuticals from arsenic or organo- arsenic compounds.	K114	Vicinals from the purification of toluenediamine in production of
K102	Residue from the use of activated carbon for decolorization in the production of veterinary pharmaceuticals from arsenic or organo-arsenic compounds.		toluenediamine via hydrogenation of dinitrotoluene.

Code	Waste description	Code	Waste description
K115	Heavy ends from purification of toluenediamine in the production of toluenediamine via hydrogenation of dinitrotoluene.	PROI CON	ARDED COMMERCIAL CHEMICAL DUCTS, OFF-SPECIFICATION SPECIES, TAINER RESIDUALS, AND SPILL DUES THEREOF-ACUTE HAZARDOUS
K116	Organic condensate from the solvent recovery column in the production of toluene diisocyanate via phosgenation of toluenediamine.	WAST	TE ALPHABETIZED LISTING CAN BE FOUND AT 40 CFR 261.33.)
K117	Wastewater from the reactor vent gas scrubber in the production of ethylene dibromide via bromination of ethene.	P001	2H-1-Benzopyran-2-one, 4-hydroxy-3-(3-oxo-1-phenylbutyl)-, & salts, when present at concentrations greater than 0.3%
K118	Spent adsorbent solids from purification of ethylene dibromide in the production of ethylene dibromide via bromination of	P001	Warfarin, & salts, when present at concentrations greater than 0.3%
	ethene.	P002	1-Acetyl-2-thiourea
K123	Process wastewater (including supernates, filtrates, and washwaters) from the production of ethylenebisdithiocarbamic acid and its salt.	P002	Acetamide, N-(aminothioxomethyl)-
		P003	2-Propenal
		P003	Acrolein
K124	Reactor vent scrubber water from the production of ethylenebisdithiocarbamic acid and its salts.	P004	1,4,5,8-Dimethanonaphthalene, 1,2,3,4,10,10-hexa-chloro-1,4,4a,5,8,8a,- hexahydro-, (1alpha, 4alpha, 4abeta, 5alpha, 8alpha, 8abeta)-
K125	Filtration, evaporation, and centrifugation solids from the production of ethylenebisdithiocarbamic acid and its salts.	P004	Aldrin
		P005	2-Propea-1-ol
		P005	Allyi alcohol
K126	Baghouse dust and floor sweepings in	P006	Aluminum phosphide (R,T)
	milling and packaging operations from production or formulation of	P007	3(2H)-Isoxazolone, 5-(aminomethyl)-
	ethylenebisdithiocarbamic acid and its	P007	5-(Aminomethyl)-3-isoxazolol
<b>1</b> /101	salts.	P008	4-Aminopyridine
K131	Wastewater from the reactor and spent sulfuric acid from the acid dryer from the	P008	4-Pyridinamine
	production of methyl bromide.	P009	Ammonium picrate (R)
K132	Spent absorbent and wastewater separator	P009	Phenol, 2,4,6-trinitro-, ammonium salt (R)
	solids from the production of methyl bromide.	P010	Arsenic acid H <sub>3</sub> AsO <sub>4</sub>
K136	Still bottoms from the purification of	P011	Arsenic oxide As <sub>2</sub> O <sub>5</sub>
	ethylene dibromide in the production of	P011	Arsenic pentoxide
	ethylene dibromide via bromination of	P012	Arsenic oxide As <sub>2</sub> O <sub>3</sub>
	ethene.	P012	Arsenic trioxide
	•	P013	Barium cyanide

Code	Waste description	Code	Waste description
P014	Benzenethiol	P036	Dichlorophenylarsine
	<del></del>	P037	2,7:3,6-Dimethanonaphth[2,3-b]oxirene,
P014	Thiophenol	2001	3,4,5,6,9,9-hexachloro-1a,2,2a,3,6,6a,7,7a-
P015	Beryllium  Bisklansmerholesher		octahydro-, (1aalpha, 2beta, 2aalpha,
P016	Dichloromethyl ether	2005	3beta, 6beta, 6aalpha, 7beta, 7aalpha)-
P016	Methane, oxybis[chloro-	P037	Dieldrin
P017	2-Propanone, 1-bromo-	P038	Arsine, diethyl-
P017	Bromoacetone	P038	Diethylarsine
P018	Brucine	P039	Disulfoton
P018	Strychnidin-10-one, 2,3-dimethoxy-	P039	Phosphorodithioic acid, O,O-diethyl S-[2-(ethylthio)ethyl] ester
P020	Dinoseb	P040	O,O-Diethyl O-pyrazinyl phosphorothioate
P020	Phenol, 2-(1-methylpropyl)-4,6-dinitro-	P040	Phosphorothioic acid, O,O-diethyl O-
P021	Calcium cyanide	PONU	pyrazinył ester
P021	Calcium cyanide Ca(CN) <sub>2</sub>	P041	Diethyl-p-nitrophenyl phosphate
P022	Carbon disulfide	P041	Phosphoric acid, diethyl 4-nitrophenyl
P023	Acetaldehyde, chloro-		ester
P023	Chloroacetaldehyde	P042	1,2-Benzenediol, 4-[1-hydroxy-2-
P024	Benzenamine, 4-chloro-		(methylamino)ethyl]-, (R)-
P024	p-Chloroaniline	P042	Epinephrine
P026	1-(o-Chlorophenyl)thiourea	P043	Diisopropyifluorophosphate (DFP)
P026	Thiourea, (2-chlorophenyi)-	P043	Phosphorofluoridic acid, bis(1- methylethyl) ester
P027	3-Chloropropionitrile	P044	Dimethoate
P027	Propanenitrile, 3-chloro-	P044	Phosphorodithioic acid, O,O-dimethyl S-
P028	Benzene, (chloromethyl)-	P044	[2-(methylamino)-2-oxoethyl] ester
P028	Benzyl chloride	P045	2-Butanone, 3,3-dimethyl-1-(methylthio)-,
P029	Copper cyanide		O-[methylamino)carbonyl] oxime
P029	Copper cyanide Cu(CN)	P045	Thiofanox
P030	Cyanides (soluble cyanide salts), not otherwise specified	P046	alpha, alpha-Dimethylphenethylamine
P031	Cyanogen	P046	Benzeneethanamine, alpha, alpha- dimethyl-
P031	Ethanedinitrile	P047	4,6-Dinitro-o-cresol, & salts
P033		P047	Phenol, 2-methyl-4,6-dinitro-, & salts
P033		P048	2,4-Dinitrophenol
P034		P048	Phenol, 2,4-dinitro-
P034		P049	Dithiobiuret
P036			

Code	Waste description	Code	Waste description
P049	Thioimidodicarbonic diamide [(H <sub>2</sub> N)C(S)] <sub>2</sub> NH	P066	Ethanimidothioic acid, N- [[(methylamino)carbonyl]oxy]-, methyl
P050	6,9-Methano-2,4,3-		ester
	benzodioxathiepin,6,7,8,9,10,10- hexachloro-1,5,5a,6,9,9a-hexahydro-,3-	P066	Methomyl
	oxide	P067	1,2-Propylenimine
P050	Endosulfan	P067	Aziridine, 2-methyl-
P051	2,7:3,6-Dimethanonaphth[2,3-b]oxirene,	P068	Hydrazine, methyl-
	3,4,5,6,9,9-hexachloro-1a,2,2a,3,6,6a,7,7a-	P068	Methyl hydrazine
	octahydro-, (1aalpha, 2beta, 2abeta, 3alpha, 6alpha, 6abeta, 7beta, 7aalpha)- &	P069	2-Methyllactonitrile
	metabolites	P069	Propanenitrile, 2-hydroxy-2-methyl-
P051	Endrin	P070	Aldicarb
2051 2054	Endrin, & metabolites Aziridine	P070	Propanal, 2-methyl-2-(methylthio)-, O- [(methylamino)carbonyl]oxime
2054	Ethyleneimine	P071	Methyl parathion
2056	Fluorine	P071	Phosphorothioic acid, O,O,-dimethyl O-(4 nitrophenyl) ester
<sup>2</sup> 0 <i>5</i> 7	Acetamide, 2-fluoro-	P072	alpha-Naphthylthiourea
057	Fluoroacetamide	P072	Thiourea, 1-naphthalenyl-
058	Acetic acid, fluoro-, sodium salt	P073	Nickel carbonyl
058	Fluoroacetic acid, sodium salt	P073	Nickel carbonyi Ni(CO)4, (T-4)-
059	4,7-Methano-1H-indene, 1,4,5,6,7,8,8-	P074	Nickel cyanide
059	heptachioro-3a,4,7,7a-tetrahydro-	P074	Nickel cyanide Ni(CN) <sub>2</sub> .
060	Heptachlor	P075	Nicotine, & salts
vou	1,4,5,8-Dimethanonaphthalene, 1,2,3,4,10,10-bexa-chloro-1,4,4a,5,8,8a,- hexahydro-, (1alpha, 4alpha, 4abeta, 5beta,	P075	Pyridine, 3-(1-methyl-2-pyrrolidinyl)-,(S)-, & salts
	Sbeta, Sabeta)-	P076	Nitric oxide
060	Iso <del>dria</del>	P076	Nitrogen oxide NO
062	Hexaethyl tetraphosphate	P077	Benzenamine, 4-nitro-
062	Tetraphosphoric acid, hexaethyl ester	P077	p-Nitroaniline
063	Hydrocyanic acid	P078	Nitrogen dioxide
063	Hydrogen cyanide	P078	Nitrogen oxide NO <sub>2</sub>
064	Methane, isocyanato-	P081	1,2,3-Propanetriol, trinitrate (R)
064	Methyl isocyanate	P081	Nitroglycerine (R)
265	Fulminic acid, mercury(2+) salt (R,T)	P082	Methanimine, N-methyl-N-nitroso-
<b>X65</b>	Mercury fulminate (R,T)	P082	N-Nitrosodimethylamine

Code	Waste description	Code	Waste description
P084	N-Nitrosomethylvinylamine	P104	Silver cyanide
2084	Vinylamine, N-methyl-N-nitroso-	P104	Silver cyanide Ag(CN)
2085	Diphosphoramide, octamethyl-	P105	Sodium azide
P085	Octamethylpyrophosphoramide	P106	Sodium cyanide
2087	Osmium oxide OsO <sub>4</sub> , (T-4)-	P106	Sodium cyanide Na(CN)
2087	Osmium tetroxide	P107	Strontium sulfide SrS
P088	7-Oxabicyclo[2.2.1]heptane-2,3-	P108	Strychnidin-10-one, & salts
	dicarboxylic acid	P108	Strychnine, & salts
P088	Endothall	P109	Tetraethyldithiopyrophosphate
P089	Parathion	P109	Thiodiphosphoric acid, tetraethyl ester
P089	Phosphorothioic acid, O,O-diethyl-O-(4-	P110	Plumbane, tetraethyl-
<b>~~~</b>	nitrophenyl) ester	P110	Tetraethyl lead
P092	Mercury, (acetato-O)phenyl-	P111	Diphosphoric acid, tetraethyl ester
P092	Phenylmercury acetate	P111	Tetraethyl pyrophosphate
P093	Phenylthiourea	P112	Methane, tetranitro- (R)
P093	Thiourea, phenyl-	P112	Tetranitromethane (R)
P094	Phorate	P113	Thallic oxide
P094	Phosphorodithioic acid, O,O-diethyl S- [(ethylthio)methyl] ester	P113	Thallium oxide Tl <sub>2</sub> O <sub>3</sub>
P095	Carbonic dichloride	P114	Selenious acid, dithallium (1+) salt
P095	Phosgene	P114	Thallium(I) selenite
P096	Hydrogen phosphide	P115	Sulfuric acid, dithallium (1+) salt
P096	Phosphine	P115	Thallium(I) sulfate
P097	Famphur	P116	Hydrazinecarbothioamide
P097	Phosphorothioic acid O-[4-	P116	Thiosemicarbazide
	[(dimethylamino)sulfonyl]phenyl] O,O-	P118	Methanethiol, trichloro-
	dimethyl ester	P118	Trichloromethanethiol
P098	Potassium cyanide	P119	Ammonium vanadate
P098	Potassium cyanide K(CN)	P119	Vanadic acid, ammonium salt
P099	Argentate (1-), bis(cyano-C)-, potassium	P120	Vanadium oxide V <sub>2</sub> O <sub>5</sub>
P099	Potassium silver cyanide	P120	Vanadium pentoxide
P101	Ethyl cyanide	P121	Zinc cyanide
P101	Propanenitrile	P121	Zinc cyanide Zn(CN) <sub>2</sub>
P102	2-Propyn-1-ol Propargyl alcohol	P122	Zinc phosphide Zn <sub>3</sub> P <sub>2</sub> , when present a concentrations greater than 10% (R,T)
P102			concentrations are afer in an item in. I

Code	Waste description	Code	Waste description
PROD CONT RESIL	ARDED COMMERCIAL CHEMICAL DUCTS, OFF-SPECIFICATION SPECIES, CAINER RESIDUES, AND SPILL DUES THEREOF-TOXIC WASTES  LPHABETIZED LISTING CAN BE	<b>U010</b>	Azirino [2',3':3,4]pyrrolo[1,2-a]indole-4,7-dione, 6-amino-8- [[(aminocarbonyl)oxy]methyl]- 1,1a,2,8,8a,8b-hexahydro-8a-methoxy-5- methyl-, [1aS-(1aalpha, 8beta, 8aalpha, 8balpha)]-
FOUN	'D AT 40 CFR 261.33.)	U010	Mitomycin C
		U011	1H-1,24-Triazol-3-amine
	2,3,4,6-Tetrachlorophenol	U011	Amitrole
	24,5-T	U012	Aniline (I,T)
	2,4,5-Trichlorophenol 2,4,6-Trichlorophenol	U012	Benzenamine (LT)
	Acetic acid, (2,4,5-trichlorophenoxy)-	U014	Auramine
See F027	Pentachlorophenol Phenol, 2,3,4,6-tetrachloro- Phenol, 2,4,5-trichloro-	U014	Benzenamine, 4,4'-carbonimidoylbis[N,N-dimethyl-
	Phenol, 2,4,6-trichloro-	U015	Azaserine
	Phenol, pentachloro-	U015	L-Serine, diazoacetate (ester)
	Propanoic acid, 2-(2,4,5- trichlorophenoxy)-	U016	Benz[c]acridine
	Silvex (2,4,5-TP)	U017	Benzal chloride
		U017	Benzene, (dichloromethyl)-
U001	Acetaldehyde (I)	U018	Benz(a)anthracene
U001	Ethanal (I)	U019	Benzene (I,T)
U002	2-Propanone (I)	U020	Benzenesulfonic acid chloride (C,R)
L'002	Acetone (I)	U020	Benzenesulfonyl chloride (C,R)
U003	Acetonitrile (I,T)	U021	[1,1'-Biphenyi]-4,4'-diamine
U004	Acetophenone	U021	Benzidine
U004	Ethanone, 1-phenyi-	U022	Benzo(a)pyrene
C005	2-Acetylaminofluorene	U023	Benzene, (trichloromethyl)-
U005	Acetamide, N-9H-fluoren-2-yl	U023	Benzotrichloride (C,R,T)
U006	Acetyl chloride (C,R,T)	U024	Dichloromethoxy ethane
U007	2-Propenamide	U024	Ethane, 1,1'-[methylenebis(oxy)]bis[2-
U007	Acrylamide	0024	chloro-
U008	2-Propenoic acid (I)	U025	Dichloroethyl ether
U008	Acrylic acid (I)	U025	Ethane, 1,1'-oxybis[2-chloro-
Ľ009	2-Propenenitrile	U026	Chlornaphazin
Ľ009	Acrylonitrile	U026	Naphthalenamine, N,N'-bis(2-chloroethyl)-

Code	Waste description	Code	Waste description
U027	Dichloroisopropyl ether	U043	Vinyl chloride
U027	Propane, 2,2'-oxybis[2-chloro-	Ú044	Chloroform
U028	1,2-Benzenedicarboxylic acid, bis(2-	U044	Methane, trichloro-
	ethylhexyl) ester	U045	Methane, chloro- (I,T)
U028	Diethylhexyl phthalate	U045	Methyl chloride (I,T)
U029	Methane, bromo-	U046	Chloromethyl methyl ether
U029	Methyl bromide	U046	Methane, chloromethoxy-
U030	4-Bromophenyl phenyl ether	U047	beta-Chloronaphthalene
U030	Benzene, 1-bromo-4-phenoxy-	U047	Naphthalene, 2-chloro-
U031	1-Butanol (I)	U048	o-Chlorophenol
U031	n-Butyi alcohol (I)	U048	Phenol, 2-chloro-
U032	Calcium chromate	U049	4-Chloro-o-toluidine, hydrochloride
U032	Chromic acid H <sub>2</sub> CrO <sub>4</sub> , calcium salt	U049	Benzenamine, 4-chloro-2-methyl-,
U033	Carbon oxyfluoride (R,T)		hydrochloride
U033	Carbonic difluoride	U050	Chrysene
U034	Acetaldehyde, trichloro-	U051	Creosote
U034	Chioral	U052	Cresol (Cresylic acid)
U035	Benzenebutanoic acid, 4-[bis(2-	U052	Phenol, methyl-
	chloroethyl)amino]-	U053	2-Butenal
U035	Chlorambucil	U053	Crotonaldehyde
U036	4,7-Methano-1H-indene, 1,2,4,5,6,7,8,8-	U055	Benzene, (1-methylethyl)- (I)
1:026	ocrachloro-2,3,3a,4,7,7a-hexahydro-	U055	Cumene (I)
U036	Chlordane, alpha & gamma isomers	U056	Benzene, hexahydro- (I)
U037	Benzene, chloro-	U0 <b>56</b>	Cyclohexane (I)
U037	Chlorobenzene	U057	Cyclohexanone (I)
U038	Benzeneacetic acid, 4-chloro-alpha-(4-chlorophenyi)-alpha-hydroxy-, ethyl ester	U058	2H-1,3,2-Oxazaphosphorin-2-amine, N,N-bis(2-chloroethyl)tetrahydro-, 2-oxide
U038	Chlorobenzilate	U058	Cyclophosphamide
Ľ039	p-Chloro-m-cresol	U059	5,12-Naphthacenedione, 8-acetyl-10-[(3-
U'039	Phenol, 4-chloro-3-methyl-		amino-2,3,6-trideoxy)-alpha-L-lyxo-
U041	Epichlorohydrin		hexopyranosyi)oxy]-7,8,9,10-tetrahydro- 6,8,11-trihydroxy-1-methoxy-, (8S-cis)-
U041	Oxirane, (chloromethyl)-	U059	Daunomycin
U'042	2-Chloroethyl vinyl ether	U060	Benzene, 1,1'-(2,2-
U042	Ethene, (2-chloroethoxy)-		dichloroethylidene)bis[4-chloro-
U043	Ethene, chloro-	U060	DDD

1.2-Dichloroethyldene   1.2-Dichloroethylene   1.2-Dichloroethylen	Code	Waste description	Code	Waste description
Unterlate   Uniterlate   Unit	Ľ'061		U079	1,2-Dichloroethylene
U062   Carbamothioic acid, bis(1-methylethyl)-, S-(2,3-dichloro-2-propenyl) ester   U080   Methylene chloride	11041		U079	Ethene, 1,2-dichloro-,(E)-
S-(2,3-dichloro-2-propenyl) ester			U080	Methane, dichloro-
Diallate	0002	S-(2.3-dichloro-2-propenyl) ester	U080	Methylene chloride
Dienza_nantracene U064 Benzo[sx]pentaphene U065 Dibenzo[a,i]pyrene U066 1,2-Dibromo-3-chloropropane U067 Ethane, 1,2-dibromo- U068 Methane, dibromo- U068 Methane, dibromo- U069 1,2-Benzenedicarboxylic acid, dibutyl ester U070 Dibutyl phthalate U070 Benzene, 1,2-dichloro- U071 Benzene, 1,3-dichloro- U072 P-Dichlorobenzene U073 1,1-Biphenyl]-4,4'-diamine, 3,3'-dichloro- U074 2-Butene, 1,4-dichloro- U075 Dichlorobenzidine U076 Ethane, di-lorodiffluoronethane U077 Biene, 1,4-dichloro- U078 Clare, 1,4-dichloro- U079 Dichlorobenzene U070 Dichlorobenzene U071 Benzene, 1,4-dichloro- U072 Dichlorobenzene U073 Dichlorobenzene U074 L-Dichloro- U075 Dichlorobenzene U076 Clare, 1,4-dichloro- U077 Clare, 1,4-dichloro- U078 Clare, 1,4-dichloro- U079 Dichlorobenzene U070 Dichlorobenzene U071 Dichlorobenzene U071 Dichlorobenzene U072 Dichlorobenzene U073 (1,1'-Biphenyl]-4,4'-diamine, 3,3'-dichloro- U074 C-Butene, 1,4-dichloro- (1,T) U075 Dichlorodiffuoromethane U076 Ethane, 1,1-dichloro- U077 Ethane, dichlorodiffuoromethane U077 Ethane, 1,2-dichloro- U078 Ethane, 1,2-dichloro- U079 Ethane, 1,2-dichloro- U070 Ethane, 1,2-dichloro- U070 Ethane, 1,2-dichloro- Ethylidene dichloride U077 Ethane, 1,2-dichloro- Ethylidene dichloride U077 Ethane, 1,1-Dichloroethylene U078 L-Dichloroethylene U079 Dimethylamine (1) U070 Ethane, 1,1-Dichloroethylene U070 Ethane, 1,1-Dichloroethylene U070 Ethane, 1,1-Dichloroethylene U071 Ethane, 1,1-Dichloroethylene U072 Dimethylaminoazobenzene U073 L-Dichloroethylene U074 L-Dichloroethylene U075 Ethane, 1,1-Dichloroethylene U076 Ethane, 1,1-Dichloroethylene U077 Ethenedichloride U078 Ethane, 1,1-Dichloroethylene U079 Ethane, 1,1-Dichloroethylene U079 Ethane, 1,1-Dichloroethylene	U062		U081	2,4-Dichlorophenol
U064   Benzo[st]pentaphene   U082   2,6-Dichlorophenol	U063	Dibenz[a,h]anthracene	U081	Phenol, 2,4-dichloro-
U064 Dibenzo[a,i]pyrene U066 1,2-Dibromo-3-chloropropane U066 Propane, 1,2-dibromo-3-chloro U067 Ethane, 1,2-dibromo- U067 Ethylene dibromide U068 Methane, dibromo- U068 Methane, dibromo- U069 1,2-Benzenedicarboxylic acid, dibutyl ester U070 Dibutyl phthalate U071 Benzene, 1,3-dichloro- U071 Benzene, 1,3-dichloro- U071 Benzene, 1,4-dichloro- U071 U071 U071 Benzene U071 U071 U071 Benzene, 1,4-dichloro- U071 U071 U071 U071 U071 U071 U071 U071	U064		U082	2,6-Dichlorophenol
1.2-Dioromo-3-chloro-  U083   Propylene dichloride	U064	• ••	U082	Phenol, 2,6-dichloro-
U066 Propane, 1,2-dibromo-3-chloro U067 Ethane, 1,2-dibromo- U068 Ethylene dibromide U068 Methane, dibromo- U069 1,2-Benzenedicarboxylic acid, dibutyl ester U069 Dibutyl phthalate U070 Benzene, 1,2-dichloro- U070 O-Dichlorobenzene U071 Benzene, 1,3-dichloro- U071 Benzene, 1,4-dichloro- U072 Benzene, 1,4-dichloro- U073 [1,1'-Biphenyl]-4,4'-diamine, 3,3'-dichloro- U074 1,4-Dichloro-2-butene (I,T) U075 Dichlorodifluoromethane U076 Ethane, 1,1-dichloro- U077 Ethane, 1,2-dichloro- U078 Ethane, 1,2-dichloro- U079 Dichlorodifluoromethylene U070 Dichlorodifluoronethylene U070 Dichlorodenide U070 Dichlorodenide U070 Dichlorodenide U070 Dichlorodenide U070 Dichlorodenide U070 Dichlorodenide U070 Dichlorodifluoro- Ethane, 1,1-dichloro- Ethylidene dichloride U070 Dimethylaminoazobenzene U071 Dimethylaminoazobenzene U071 Dichlorocthylene U071 Dichlorothylene U071 Dimethylaminoazobenzene U071 Dimethylaminoazobenzene U071 Dichlorothylene U071 Dichlorothylene U072 Dimethylaminoazobenzene U073 Dimethylaminoazobenzene U074 Dimethylaminoazobenzene U075 Dimethylaminoazobenzene U076 Dimethylaminoazobenzene U077 Dimethylaminoazobenzene	U066	* ***	U083	Propane, 1,2-dichloro-
U067 Ethylene dibromo- U068 Methane, dibromo- U068 Methylene bromide U069 1,2-Benzenedicarboxylic acid, dibutyl ester U069 Dibutyl phthalate U0700 Benzene, 1,2-dichloro- U0701 Benzene, 1,3-dichloro- U0701 Benzene, 1,3-dichloro- U0702 Benzene, 1,4-dichloro- U0703 [1,1'-Biphenyl]-4,4'-diamine, 3,3'-dichloro- U0704 1,4-Dichloro-2-butene (I,T) U0705 Dichlorodifluoromethane U0706 Ethylene dichloride U0707 Ethylene dichloride U0708 Ethylipene (I,1-dichloro- U0709 Dichlorodifluoro- U0709 Dichlorodifluoro- U0709 Dichlorodifluoro- U0709 Dichlorodifluoro- Ethylene dichloride U0709 Ethylene dichloride U0709 Ethylene dichloride U0709 Ethylene dichloride U0709 Ethylipene (I,1-dichloro- Ethylene dichloride U0709 Ethylene dichloride U0709 Ethylene dichloride U0709 Ethylene dichloride U0709 Dimethylamino (N-dimethyl-4- (phenylazo)- U0709 Dimethylaminoazobenzene U0710 Dimethylaminoazobenzene U0710 Dimethylaminoazobenzene U0710 Dimethylaminoazobenzene U0710 Dimethylaminoazobenzene	U066		U083	Propylene dichloride
U068 Methane, dibromo- U069 Methylene bromide U069 1,2-Benzenedicarboxylic acid, dibutyl ester U060 Dibutyl phthalate U070 Benzene, 1,2-dichloro- U070 o-Dichlorobenzene U071 Benzene, 1,3-dichloro- U071 Benzene, 1,4-dichloro- U071 Benzene, 1,4-dichloro- U072 Benzene, 1,4-dichloro- U073 [1,1'-Biphenyl]-4,4'-diamine, 3,3'-dichloro- U073 3,3'-Dichlorobenzidine U074 2-Butene, 1,4-dichloro- U075 Dichlorodenizidine U076 Ethane, 1,1-dichloro- U077 Ethane, 1,2-dichloro- U078 Ethane, 1,2-dichloro- U079 Ethane, 1,2-dichloro- U070 Dimethylamine (I) U070 Dimethylamine, 3,3'-Dimethoxybenzidine U071 Dimethylamine, 3,3'-Dimethylamine,	U067		U084	1,3-Dichloropropene
U068 Methane, dibromo- U069 Methylene bromide U069 1.2-Benzenedicarboxylic acid, dibutyl ester U069 Dibutyl phthalate U070 Benzene, 1,2-dichloro- U070 o-Dichlorobenzene U071 Benzene, 1,3-dichloro- U071 Benzene, 1,4-dichloro- U071 Benzene, 1,4-dichloro- U071 Benzene, 1,4-dichloro- U072 P-Dichlorobenzene U073 J3'-Dichlorobenzene U073 J3'-Dichlorobenzene U074 L4-Dichloro-2-butene (I,T) U075 Dichlorodifluoromethane U076 Ethane, 1,1-dichloro- U077 Ethane, 1,2-dichloro- U077 Ethane, 1,2-dichloro- Ethylene dichloride U077 Ethane, 1,2-dichloro- Ethylene dichloride U078 L2-Benzenedicarboxylic acid, diethyl ester U089 Diethylstilbesterol U089 Phenol, 4,4'-(1,2-diethyl-1,2-ethenediyl)bis, (E)- U090 Dihydrosafrole U091 J3-Benzodioxole, 5-propyl- U092 Dimethylamine, 3,3'-dimethoxy- U093 J3'-Dimethoxybenzidine U094 L3-Binzendioxole U095 Dimethylamine (I) U096 Ethane, 1,1-dichloro- Ethylidene dichloride U097 Ethane, 1,2-dichloro- Ethylene dichloride U098 Phenol, 4,4'-diamine, 3,3'-dimethyl-4- (phenylazo)- P-Dimethylaminoazobenzene U099 Dimethylaminoazobenzene U090 Dimethylaminoazobenzene U091 J3-Dimethylaminoazobenzene U092 Dimethylaminoazobenzene	U067	Ethylene dibromide	U084	1-Propene, 1,3-dichloro-
U069 1,2-Benzenedicarboxylic acid, dibutyl ester U069 Dibutyl phthalate U070 Benzene, 1,2-dichloro- U070 o-Dichlorobenzene U071 Benzene, 1,3-dichloro- U071 m-Dichlorobenzene U072 Benzene, 1,4-dichloro- U073 [1,1'-Biphenyl]-4,4'-diamine, 3,3'-dichloro- U073 3,3'-Dichlorobenzidine U074 1,4-Dichloro-2-butene (I,T) U075 Dichlorodifluoromethane U076 Ethane, 1,1-dichloro- U077 Ethane, 1,2-dichloro- U078 Ethane, 1,2-dichloro- U079 Ethane, 1,2-dichloro- U070 Ethane, 1,2-dichloro- U070 Ethane, 1,2-dichloro- U070 Ethane, 1,2-dichloro- U070 Ethylene dichloride U071 Ethane, 1,2-dichloro- U071 Ethane, 1,2-dichloro- U071 Ethane, 1,1-dichloro- U071 Ethane, 1,2-dichloro- U071 Ethane, 1,1-Dichloroethylene U071 Ethane, 1,1-Dichloroethylene U071 Injentylaminoazobenzene U071 Injentylaminoazobenzene U072 Injentylaminoazobenzene U073 Injentylaminoazobenzene U074 Injentylaminoazobenzene U075 Injentylaminoazobenzene U076 Ethane, 1,2-dichloro- U077 Ethane, 1,2-dichloro- U077 Ethane, 1,2-dichloro- U078 Injentylaminoazobenzene U079 Injentylaminoazobenzene U070 Injentylaminoazobenzene U071 Injentylaminoazobenzene U071 Injentylaminoazobenzene U072 Injentylaminoazobenzene	U068		U085	1,2:3,4-Diepoxybutane (I,T)
U069 1,2-Benzenedicarboxylic acid, dibutyl ester U069 Dibutyl phthalate U070 Benzene, 1,2-dichloro- U070 o-Dichlorobenzene U071 Benzene, 1,3-dichloro- U071 Benzene, 1,3-dichloro- U072 Benzene, 1,4-dichloro- U073 [1,1'-Biphenyl]-4,4'-diamine, 3,3'-dichloro- U074 1,4-Dichloro-2-butene (I,T) U075 Dichlorodiffuoromethane U076 Ethane, 1,1-dichloro- U077 Ethane, 1,2-dichloro- U078 Ethane, 1,2-dichloro- U079 Ethane, 1,2-dichloro- U070 Ethane, 1,2-dichloro- U070 Ethane, 1,2-dichloro- U071 Ethane, 1,2-dichloro- U071 Ethane, 1,2-dichloro- U072 Ethane, 1,2-dichloro- U073 Ethane, 1,2-dichloro- U074 Ethane, 1,2-dichloro- U075 Ethane, 1,2-dichloro- U076 Ethylidene dichloride U077 Ethane, 1,2-dichloro- U078 Ethane, 1,2-dichloro- U079 Ethane, 1,2-dichloro- U070 Ethane, 1,2-dichloro- U071 Ethane, 1,2-dichloro- U071 Ethane, 1,2-dichloro- U072 Ethane, 1,2-dichloro- U073 L1-Dichloroethylene U074 L1-Dichloroethylene U075 L2 Dimethylaminoazobenzene U076 L2 Dimethylaminoazobenzene U077 L2 Dimethylaminoazobenzene U078 L2 Dimethylaminoazobenzene U079 L2 Dimethylaminoazobenzene	U068	Methylene bromide	U <b>085</b>	2,2'-Bioxirane
U069 Dibutyl phthalate U070 Benzene, 1,2-dichloro- U070 o-Dichlorobenzene U071 Benzene, 1,3-dichloro- U071 m-Dichlorobenzene U072 Benzene, 1,4-dichloro- U073 [1,1'-Biphenyl]-4,4'-diamine, 3,3'-dichloro- U074 1,4-Dichloro-2-butene (I,T) U075 Dichlorodifluoro- U076 Ethane, 1,1-dichloro- U077 Ethane, 1,2-dichloro- U078 Ethyliene dichloride U079 Ethane, 1,2-dichloro- U070 Dimethylamine, N,N-dimethyl-4- U070 Ethylichoroethylene U080 Diethylstilbesterol U089 Diethylstilbesterol U089 Phenol, 4,4'-(1,2-diethyl-1,2-ethenediyl)bis, (E)- U070 1,3-Benzodioxole, 5-propyl- U170 1,1-Biphenyl]-4,4'-diamine, 3,3'-dimethoxy- U079 Dichlorodifluoro- U070 Ethane, 1,1-dichloro- U070 Ethane, 1,2-dichloro- U070 Ethylidene dichloride U070 Ethylidene dichloride U070 CO71 Interhylamine, N-methyl-(I) U071 Ethylene dichloride U072 Dimethylaminoazobenzene U073 Intertylaminoazobenzene U074 Interhylaminoazobenzene U075 Dimethylaminoazobenzene U076 Ethylene dichloride U077 Ethylene dichloride U078 Interhylaminoazobenzene U079 Dimethylaminoazobenzene	U069	-	U086	Hydrazine, 1,2-diethyl-
U070 Benzene, 1,2-dichloro- U070 o-Dichlorobenzene U071 Benzene, 1,3-dichloro- U071 m-Dichlorobenzene U072 Benzene, 1,4-dichloro- U073 p-Dichlorobenzene U074 [1,1'-Biphenyl]-4,4'-diamine, 3,3'-dichloro- U075 [1,4-Dichloro-2-butene (I,T) U076 Dichlorodiffuoromethane U077 Dichlorodiffuoro- U078 Dichlorodiffuoro- U079 Dichlorodiffuoro- U079 Dichlorodiffuoro- U070 Dichlorodiffuoro- U070 Dichlorodiffuoro- U071 Dichlorodiffuoro- U071 Dichlorodiffuoro- U072 Dimethylamine, 3,3'-Dimethoxybenzidine U073 Dichlorodiffuoro- U074 Dichlorodiffuoro- U075 Dichlorodiffuoro- U076 Ethane, 1,1-dichloro- U077 Ethane, 1,2-dichloro- U077 Ethylidene dichloride U077 Dimethylamine, N-methyl-(I) U077 Ethylidene dichloride U078 Dimethylamine, N-methyl-(I) U079 Ethylidene dichloride U070 Dimethylaminoazobenzene U071 Dimethylaminoazobenzene U071 Dimethylaminoazobenzene U071 Dimethylaminoazobenzene U071 Dimethylaminoazobenzene	U069		U086	N,N'-Diethylhydrazine
U070 o-Dichlorobenzene U071 Benzene, 1,3-dichloro- U071 m-Dichlorobenzene U072 Benzene, 1,4-dichloro- U073 p-Dichlorobenzene U075 [1,1'-Biphenyl]-4,4'-diamine, 3,3'-dichloro- U076 [1,4-Dichloro-2-butene (I,T) U077 Dichlorodiffuoromethane U078 Methane, dichloro- U079 Ethane, 1,1-dichloro- U070 Ethane, 1,2-dichloro- U070 Ethylidene dichloride U070 Ethane, 1,2-dichloro- U071 Ethane, 1,2-dichloro- U072 Ethylidene dichloride U073 Ethylene dichloride U074 Ethane, 1,2-dichloro- U075 Ethylidene dichloride U076 Ethane, 1,2-dichloro- U077 Ethylene dichloride U077 Ethylene dichloride U078 I,1-Dichloroethylene U079 Dimethylaminoazobenzene U070 Dimethylaminoazobenzene U071 Dimethylaminoazobenzene U071 Dimethylaminoazobenzene U072 Dimethylaminoazobenzene	U070		U <b>087</b>	O,O-Diethyl S-methyl dithiophosphate
U071 m-Dichlorobenzene U072 Benzene, 1,4-dichloro- U073 Benzene, 1,4-dichloro- U074 p-Dichlorobenzene U075 [1,1'-Biphenyl]-4,4'-diamine, 3,3'-dichloro- U076 1,4-Dichloro-2-butene (I,T) U077 Dichlorodifluoromethane U078 Methane, dichloro- U079 Ethane, 1,1-dichloro- U070 Ethylidene dichloride U070 Ethylidene dichloride U071 Ethylidene dichloride U071 I,2-Benzenedicarboxylic acid, diethyl ester U089 Diethylstilbesterol U089 Phenol, 4,4'-(1,2-diethyl-1,2-ethenediyl)bis, (E)- U090 I,3-Benzodioxole, 5-propyl- U090 Dihydrosafrole U091 [1,1'-Biphenyl]-4,4'-diamine, 3,3'-dimethoxy- dimethoxy- U091 [1,1'-Biphenyl]-4,4'-diamine, 3,3'-dimethoxy- U092 Dimethylamine (I) U093 Benzenamine, N-methyl- (I) U094 Ethylidene dichloride U095 Benzenamine, N,N-dimethyl-4- (phenylazo)- U096 Phenol, 4,4'-(1,2-diethyl-1,2-ethenediyl)bis, (E)- U090 I,3-Benzodioxole, 5-propyl- U091 [1,1'-Biphenyl]-4,4'-diamine, 3,3'- dimethoxy- U091 Benzenamine, N,N-dimethyl-4- (phenylazo)- U092 Dimethylaminoazobenzene U093 Benzenamine, N,N-dimethyl-4- (phenylazo)- U096 Phenol, 4,4'-(1,2-diethyl-1,2-ethenediyl)bis, (E)- U099 Phenol, 4,4'-(1,2-diethyl-1,2-ethenediyl)bis, (E)- U090 Dihydrosafrole U090 I,3-Benzodioxole, 5-propyl- U091 [1,1'-Biphenyl]-4,4'-diamine, 3,3'- dimethoxy- dimethoxy- U091 Benzene, 1,4-dichloro- U092 Dimethylamine (I) U093 Benzenamine, N,N-dimethyl-4- (phenylazo)- U094 I,4-Dichloro- U095 Phenol, 4,4'-(1,2-diethyl-1,2-ethenediyl)bis, (E)- U096 Phenol, 4,4'-(1,2-diethyl-1,2-ethenediyl)bis, (E)- U097 Dihydrosafrole U098 Phenol, 4,4'-(1,2-diethyl-1,2-ethenediyl)bis, (E)- U099 Dihydrosafrole U099 I,3-Benzodioxole, 5-propyl- U090 Dihydrosafrole U090 I,3-Benzodioxole, 5-propyl- U090 Dihydrosafrole U091 I,3-Benzodioxole, 5-propyl- U091 I,3-Benzodioxole	C070		U087	Phosphorodithioic acid, O,O-diethyl S-
U071 m-Dichlorobenzene U072 Benzene, 1,4-dichloro- U073 p-Dichlorobenzene U073 [1,1'-Biphenyl]-4,4'-diamine, 3,3'-dichloro- U074 1,4-Dichloro-2-butene (I,T) U075 Dichlorodifluoromethane U076 Ethane, 1,1-dichloro- U076 Ethane, 1,2-dichloro- U077 Ethane, 1,2-dichloro- U078 [1,1-Dichlorobenzene U089 Phenol, 4,4'-(1,2-diethyl-1,2-ethenediyl)bis, (E)- U090 1,3-Benzodioxole, 5-propyl- U090 Dihydrosafrole U091 [1,1'-Biphenyl]-4,4'-diamine, 3,3'-dimethoxy- U095 Dichlorodifluoro- U096 Ethane, 1,1-dichloro- U097 Dimethylamine (I) U097 Ethane, 1,2-dichloro- U098 Diethylstilbesterol U089 Phenol, 4,4'-(1,2-diethyl-1,2-ethenediyl)bis, (E)- U090 1,3-Benzodioxole, 5-propyl- U091 [1,1'-Biphenyl]-4,4'-diamine, 3,3'-dimethoxy- U091 3,3'-Dimethoxybenzidine U092 Dimethylamine (I) U093 Benzenamine, N-methyl- (I) U094 Ethane, 1,2-dichloro- U095 Benzenamine, N,N-dimethyl-4- (phenylazo)- U096 Inthylaminoazobenzene U097 P-Dimethylaminoazobenzene	U071	Benzene, 1,3-dichloro-	T 1000	•
U072 Benzene, 1,4-dichloro- U073 p-Dichlorobenzene U073 [1,1'-Biphenyl]-4,4'-diamine, 3,3'-dichloro- U073 3,3'-Dichlorobenzidine U074 1,4-Dichloro-2-butene (I,T) U075 Dichlorodifluoromethane U076 Ethane, 1,1-dichloro- U076 Ethylidene dichloride U077 Ethane, 1,2-dichloro- U078 1,1-Dichloro- U079 Diethylstilbesterol U089 Phenol, 4,4'-(1,2-diethyl-1,2-ethenediyl)bis, (E)- U090 1,3-Benzodioxole, 5-propyl- U090 Dihydrosafrole U091 [1,1'-Biphenyl]-4,4'-diamine, 3,3'-dimethoxy- U091 3,3'-Dimethoxybenzidine U092 Dimethylamine (I) U093 Benzenamine, N-methyl- (I) U094 Ethane, 1,2-dichloro- U095 Benzenamine, N,N-dimethyl-4-phenylazo)- U096 I,1-Dichloroethylene	U071	m-Dichlorobenzene		
U072 p-Dichlorobenzene U073 [1,1'-Biphenyl]-4,4'-diamine, 3,3'-dichloro- U073 3,3'-Dichlorobenzidine U074 1,4-Dichloro-2-butene (I,T) U075 Dichlorodifluoromethane U075 Methane, dichlorodifluoro- U076 Ethane, 1,1-dichloro- U076 Ethylidene dichloride U077 Ethylene dichloride U078 1,1-Dichloroethylene U079 Phenol, 4,4'-(1,2-diethyl-1,2-ethenediyl)bis, (E)- U090 1,3-Benzodioxole, 5-propyl- U090 Dihydrosafrole U091 [1,1'-Biphenyl]-4,4'-diamine, 3,3'-dimethoxy- U091 [1,1'-Biphenyl]-4,4'-diamine, 3,3'-dimethoxy- U091 3,3'-Dimethoxybenzidine U092 Dimethylamine (I) U093 Benzenamine, N-methyl- (I) U094 U095 Hethanamine, N-methyl- (I) U097 Ethylene dichloride U098 Phenol, 4,4'-(1,2-diethyl-1,2-ethenediyl)bis, (E)- U090 Dihydrosafrole U091 [1,1'-Biphenyl]-4,4'-diamine, 3,3'-dimethoxy- U091 [1,1'-Biphenyl]-4,4'-diamine, 3,3'-dimethoxy- U091 3,3'-Dimethoxybenzidine U092 Dimethylamine (I) U093 Benzenamine, N-methyl- (I) U094 U095 Phenol, 4,4'-(1,2-diethyl-1,2-ethenediyl)bis, (E)- U096 Phenol, 4,4'-(1,2-diethyl-1,2-ethenediyl)bis, (E)- U096 Phenol, 4,4'-(1,2-diethyl-1,2-ethenediyl)bis, (E)- U097 U096 Phenol, 4,4'-(1,2-diethyl-1,2-ethyl-1,2-ethyl-1,2-ethyl-1,2-ethyl-1,2-ethyl-1,2-ethyl-1,2-ethyl-1,2-ethyl-1,2-ethyl-1,2-ethyl-1,2-ethyl-1,2-ethyl-1,2-e	L'072	Benzene, 1,4-dichloro-		• •
U073 [1,1'-Biphenyl]-4,4'-diamine, 3,3'-dichloro- U073 3,3'-Dichlorobenzidine U090 1,3-Benzodioxole, 5-propyi- U074 1,4-Dichloro-2-butene (I,T) U090 Dihydrosafrole U074 2-Butene, 1,4-dichloro- (I,T) U091 [1,1'-Biphenyl]-4,4'-diamine, 3,3'- U075 Dichlorodifluoromethane U091 3,3'-Dimethoxybenzidine U076 Ethane, 1,1-dichloro- U092 Dimethylamine (I) U076 Ethylidene dichloride U092 Methanamine, N-methyl- (I) U077 Ethane, 1,2-dichloro- U093 Benzenamine, N,N-dimethyl-4- U077 Ethylene dichloride U078 1,1-Dichloroethylene	U072	p-Dichlorobenzene		•
U073 3,3'-Dichlorobenzidine U090 1,3-Benzodioxole, 5-propyl- U074 1,4-Dichloro-2-butene (I,T) U090 Dihydrosafrole U091 [1,1'-Biphenyl]-4,4'-diamine, 3,3'- U095 Dichlorodifluoromethane U096 Ethane, dichlorodifluoro- U097 Methane, dichlorodifluoro- U098 Dimethylamine (I) U099 Dimethylamine (I) U099 Dimethylamine (I) U099 Dimethylamine, N-methyl- (I) U099 Ethane, 1,2-dichloro- U099 Benzenamine, N,N-dimethyl-4- (phenylazo)- U099 Dimethylaminoazobenzene U099 Dimethylaminoazobenzene	U073	[1,1'-Biphenyl]-4,4'-diamine, 3,3'-dichloro-	0069	(E)-
U074 1,4-Dichloro-2-butene (I,T) U090 Dihydrosafrole U074 2-Butene, 1,4-dichloro- (I,T) U091 [1,1'-Biphenyl]-4,4'-diamine, 3,3'- dimethoxy- U075 Dichlorodifluoromethane U076 Ethane, 1,1-dichloro- U076 Ethylidene dichloride U077 Ethane, 1,2-dichloro- U077 Ethane, 1,2-dichloro- U077 Ethylene dichloride U078 1,1-Dichloroethylene U079 Dimethylamine (I) U070 Ethylidene dichloride U091 Benzenamine, N-methyl- (I) U071 Ethylene dichloride U092 Dimethylamine, N-methyl- (I) U073 Benzenamine, N,N-dimethyl-4- (phenylazo)- U074 T12 Dimethylaminoazobenzene	U073	3,3'-Dichlorobenzidine	U090	
U074 2-Butene, 1,4-dichloro- (I,T) U075 Dichlorodifluoromethane U076 Ethane, 1,1-dichloro- U076 Ethylidene dichloride U077 Ethane, 1,2-dichloro- U077 Ethane, 1,2-dichloro- U077 Ethylene dichloride U078 1,1-Dichloroethylene U078 1,1-Dichloroethylene	U074	1,4-Dichloro-2-butene (I,T)	U090	• ••
U075 Dichlorodifluoromethane U075 Methane, dichlorodifluoro- U076 Ethane, 1,1-dichloro- U076 Ethylidene dichloride U077 Ethane, 1,2-dichloro- U077 Ethane, 1,2-dichloro- U077 Ethylene dichloride U077 Ethylene dichloride U077 Ethylene dichloride U078 1,1-Dichloroethylene U078 1,1-Dichloroethylene U079 Dimethylamine (I) U070 Benzenamine, N,N-dimethyl-4- (phenylazo)- U071 p-Dimethylaminoazobenzene	U074	2-Butene, 1,4-dichloro- (I,T)	U091	•
U076 Ethane, 1,1-dichloro- U077 Ethane, 1,2-dichloro- U077 Ethane, 1,2-dichloro- U077 Ethylene dichloride U077 Ethylene dichloride U077 Ethylene dichloride U078 1,1-Dichloroethylene U078 1,1-Dichloroethylene U079 Dimethylamine (I) U092 Methanamine, N-methyl- (I) U093 Benzenamine, N,N-dimethyl-4- (phenylazo)- U093 p-Dimethylaminoazobenzene	Ľ075	Dichlorodiffuoromethane		
U076 Ethylidene dichloride U077 Ethane, 1,2-dichloro- U077 Ethylene dichloride U077 Ethylene dichloride U078 1,1-Dichloroethylene U078 T12 Dimethylaminoazobenzene	U075	Methane, dichlorodifluoro-	U091	3,3'-Dimethoxybenzidine
U077 Ethane, 1,2-dichloro- U077 Ethylene dichloride U078 1,1-Dichloroethylene  U078 Table 1,1-Dichloroethylene  U079 Benzenamine, N,N-dimethyl-4- (phenylazo)-  U079 p-Dimethylaminoazobenzene	J076	Ethane, 1,1-dichloro-	L'092	Dimethylamine (I)
U077 Ethylene dichloride  U078 1,1-Dichloroethylene  (phenylazo)-  p-Dimethylaminoazobenzene	J076	Ethylidene dichloride	U092	Methanamine, N-methyl- (I)
C078 1,1-Dichloroethylene C078 p-Dimethylaminoazobenzene		Ethane, 1,2-dichloro-	U093	
1,1-Dichioroethylene	J077	Ethylene dichloride	1	
U078 Ethene, 1,1-dichloro-		1,1-Dichloroethylene		
	2078	Ethene, 1,1-dichloro-	C UM	., Lz-Dimetnyloenz(ajanthracene

Code	Waste description	Code	Waste description
U094	Benz[a]anthracene, 7,12-dimethyl-	U112	Ethyl acetate (I)
U095	[1,1'-Biphenyi]-4,4'-diamine, 3,3'-dimethyl-	U113	2-Propenoic acid, ethyl ester (I)
U095	3,3'-Dimethylbenzidine	U113	Ethyl acrylate (I)
U096	alpha,alpha-Dimethylbenzylhydroperoxide (R)	U114	Carbamodithioic acid, 1,2-ethanediylbis-, salts & esters
U096	Hydroperoxide, 1-methyl-1-phenylethyl-(R)	U114	Ethylenebisdithiocarbamic acid, salts & esters
U097	Carbamic chloride, dimethyl-	U115	Ethylene oxide (I,T)
U097	Dimethylcarbamoyl chloride	U115	Oxirane (I,T)
U098	1,1-Dimethylhydrazine	U116	2-Imidazolidinethione
U098	Hydrazine, 1,1-dimethyl-	U116	Ethylenethiourea
U099	1,2-Dimethylhydrazine	U117	Ethane, 1,1'-oxybis-(I)
U099	Hydrazine, 1,2-dimethyl-	U117	Ethyl ether (I)
U101	2,4-Dimethylphenol	U118	2-Propenoic acid, 2-methyl-, ethyl ester
U101	Phenol, 2,4-dimethyl-	U118	Ethyl methacrylate
U102	1,2-Benzenedicarboxylic acid, dimethyl	U119	Ethyl methanesulfonate
	ester	U119	Methanesulfonic acid, ethyl ester
U102	Dimethyl phthalate	U120	Fluoranthene
U103	Dimethyl sulfate	U121	Methane, trichlorofluoro-
U103	Sulfuric acid, dimethyl ester	U121	Trichloromonofluoromethane
U105	2,4-Dinitrotoluene	U122	Formaldehyde
U105	Benzene, 1-methyl-2,4-dinitro-	U123	Formic acid (C,T)
U106	2,6-Dinitrotoluene	U124	Furan (I)
U106	Benzene, 2-methyl-1,3-dinitro-	U124	Furfuras (I)
U107	1,2-Benzenedicarboxylic acid, dioctyl ester	U125	2-Furancarboxaldehyde (I)
U107	Di-n-octyl phthalate	U125	Furfural (I)
U108	1,4-Diethyleneoxide	U126	Glycidylaldehyde
U108	1,4-Dioxans	U126	Oxiranecarboxyaldehyde.
U109	1,2-Diphenylhydrazine	U127	Benzene, bexachioro-
U109	Hydrazine, 1,2-diphenyl-	U127	Hexachlorobenzene
U110	1-Propanimine, N-propyl-(I)	U128	1,3-Butadiene, 1,1,2,3,4,4-hexachloro-
U110	Dipropylamine (I)	U128	Hexachlorobutadiene
U111	1-Propanamine, N-nitroso-N-propyl-	U129	Cyclohexane, 1,2,3,4,5,6-hexachloro-,
U111	Di-n-propylnitrosamine		(1alpha, 2alpha, 3beta, 4alpha, 5alpha,
U112	Acetic acid ethyl ester (I)		6beta)-

Code	Waste description	Code	Waste description
11120	, The James	•••	
U129	Lindane	U146	Lead subacetate
U130	1,3-Cyclopentadiene, 1,2,3,4,5,5-, hexachloro-	U146	Lead, bis(acetato-O)tetrahydroxytri-
U130	Hexachlorocyc: pentadiene	U147	2,5-Furandione
U131	Ethane, hexachi ro-	U147	Maleic anhydride
U131	Hexachloroethan	U148	3,6-Pyridazinedione, 1.2 dihydro-
U132	Hexachlorophene	U148	Maleic hydrazide
U132	Phenol, 2,2'-methana sbis[3,4,6-trichloro-	U149	Malononitrile
U133	Hydrazine (R,T)	U149	Propanedinitrile
U134	Hydrofluoric acid ( IT)	U150	L-Phenylalanine, 4-[bis(2-chloroethyl)amino]-
U134	Hydrogen fluoride (C,T)	U150	Melphalan
U135	Hydrogen sulfide	U151	Mercury
U135	Hydrogen suifide H <sub>2</sub> S	U152	2-Propenenitrile, 2-methyl- (I,T)
U136	Arsinic acid, dimethyl-	U152	Methacrylonitrile (I,T)
U136	Cacodylic acid	U153	Methanethiol (I,T)
U137	Indeno[1,2,3-cd]pyrene	U153	Thiomethanol (I,T)
U138	Methane, iodo-	U154	Methanol (I)
U 138	Methyl iodide	U154	Methyl alcohol (I)
U140	1-Propanol, 2-methyl- (I,T)	U155	1,2-Ethanediamine, N,N-dimethyl-N'-2-
U140	Isobutyl alcohol (I,T)		pyridinyl-N'-(2-thienylmethyl)-
U141	1,3-Benzodioxole, 5-(1-propenyl)-	U155	Methapyrilene
U141	Isosafrole	U156	Carbonochloridic acid, methyl ester, (I,T)
U142	1,3,4-Metheno-2H-cyclobuta[cd]pentalen-	U156	Methyl chlorocarbonate (I,T)
	2-one, 1,1a,3,3a,4,5,5,5a,5b,6-	U157	3-Methylcholanthrene
U142	decachlorooctahydro- Kepome	U157	Benz[j]aceanthrylene, 1,2-dihydro-3-methyl-
U143	2-Butenoic acid, 2-methyl-, 7-[[2,3-	U158	4,4'-Methylenebis(2-chloroaniline)
	dihydroxy-2-(1-methoxyethyl)-3-methyl-1-	U158	Benzenamine, 4,4'-methylenebis(2-chloro-
	oxobutoxy]methyl]-2,3,5,7a-tetrahydro-1H-pyrrolizin-1-yl ester, [1S-[1alpha(Z),	U159	2-Butanone (I,T)
	7(2S°,3R°), 7aalpha]]-	U159	Methyl ethyl ketone (MEK) (I,T)
U143	Lasiocarpine	U160	2-Butanone, peroxide (R,T)
U144	Acetic acid, lead(2+) salt	U160	Methyl ethyl ketone peroxide (R,T)
J144	Lead acetate	U161	4-Methyl-2-pentanone (I)
U145	Lead phosphate	U161	Methyl isobutyl ketone (I)
J145	Phosphoric acid, lead(2+) salt (2:3)	U161	Pentanoi, 4-methyl-

Code	Waste description	Code	Waste description
U162	2-Propenoic acid, 2-methyl-, methyl ester	U180	N-Nitrosopyrrolidine
	(I,T)	U180	Pyrrolidine, 1-nitroso-
U162	Methyl methacrylate (I,T)	U181	5-Nitro-o-toluidine
U163	Guanidine, N-methyl-N'-nitro-N-nitroso-	U181	Benzenamine, 2-methyl-5-nitro
U163	MNNG	U182	1,3,5-Trioxane, 2,4,6-trimethyl-
U164	4(1H)-Pyrimidinone, 2,3-dihydro-6- methyl-2-thioxo-	U182	Paraldehyde
U164	Methylthiouracil	U183	Benzene, pentachloro-
U165	Naphthalene	U183	Pentachlorobenzene
U166	1,4-Naphthalenedione	U184	Ethane, pentachloro-
U166	•	U184	Pentachloroethane
	1,4-Naphthoquinone	U185	Benzene, pentachloronitro-
U167	1-Napthalenamine	U185	Pentachloronitrobenzene (PCNB)
U167	alpha-Naphthylamine	U186	1,3-Pentadiene (I)
U168	2-Napthalenamine	U186	1-Methylbutadiene (I)
U168	beta-Naphthylamine	U187	Acetamide, N-(4-ethoxyphenyl)-
U169	Benzene, nitro-	U187	Phenacetin
U169	Nitrobenzene (I,T)	U188	Phenol
U170	p-Nitrophenol	U189	Phosphorus suifide (R)
U170	Phenol, 4-nitro-	U189	Sulfur phosphide (R)
U171	2-Nitropropane (I,T)	U190	1,3-Isobenzofurandione
U171	Propane, 2-nitro- (I,T)	U190	Phthalic anhydride
U172	1-Butanamine, N-butyl-N-nitroso-	U191	2-Picoline
U172	N-Nitrosodi-n-butylamine	U191	Pyridine, 2-methyl-
U173	Ethanol, 2,2'-(nitrosoimino)bis-	U192	Benzamide, 3,5-dichloro-N-(1,1-dimethyl-
U173	N-Nitrosodiethanolamine		2-propynyl)-
U174	Ethanamine, N-ethyl-N-nitroso-	U192	Pronamide
U174	N-Nitrosodiethylamine	U193	1,2-Oxathiolane, 2,2-dioxide
U176	N-Nitroso-N-ethylurea	U193	1,3-Propane sultone
U176	Urea, N-ethyl-N-nitroso-	U194	1-Propanamine (I,T)
U177	N-Nitroso-N-methylurea	U194	n-Propylamine (I,T)
U177	Urea, N-methyl-N-nitroso-	U196	Pyridine
U178	Carbamic acid, methylnitroso-, ethyl ester	U197	2,5-Cyclobexadiene-1,4-dione
U178	N-Nitroso-N-methylurethane	U197	p-Benzoquinone
U179	N-Nitrosopiperidine	U200	Reserpine
U179	Piperidine, 1-nitroso-		

Code	Waste description	Code	Waste description
U200	Yohimban-16-carboxylic acid, 11,17-dimethoxy-18-[(3,4,5-trimethoxybenzoyl)oxy]-, methyl ester, (3beta, 16beta, 17alpha, 18beta, 20alpha)-	U216	Thailium chloride Tlcl
		U216	Thallium(I) chloride
		U217	Nitric acid, thallium(1+) salt
U201	1,3-Benzenediol	U217	Thallium(I) nitrate
U201	Resorcinol	U218	Ethanethioamide
U202	1,2-Benzisothiazol-3(2H)-one, 1,1-dioxide,	U218	Thioacetamide
	& salts	U219	Thiourea
U202	Saccharin, & salts	U220	Benzene, methyi-
U203	1,3-Benzodioxole, 5-(2-propenyl)-	U220	Toluene
U203	Safrole	U221	Benzenediamine, ar-methyl-
U204	Selenious acid	U221	Toluenediamine
U204	Selenium dioxide	U222	Benzenamine, 2-methyl-, hydrochloride
U205	Selenium sulfide	U222	o-Toluidine hydrochloride
U205	Selenium sulfide SeS <sub>2</sub> (R,T)	U223	Benzene, 1,3-diisocyanatomethyl- (R,T)
U206	D-Glucose, 2-deoxy-2-	U223	Toluene diisocyanate (R,T)
T 120¢	[[(methylnitrosoamino)-carbonyl]amino]-	U225	Bromoform
U206	Glucopyranose, 2-deoxy-2-(3-methyl-3-nitrosoureido)-,D-	U225	Methane, tribromo-
U206	Streptozotocin	U226	Ethane, 1,1,1-trichloro-
U207	1,2,4,5-Tetrachlorobenzene	U226	Methyl chloroform
U207	Benzene, 1,2,4,5-tetrachloro-	U227	1,1,2-Trichloroethane
U208	1,1,1,2-Tetrachloroethane	U227	Ethane, 1,1,2-trichloro-
U208	Ethane, 1,1,1,2-tetrachloro-	U228	Ethene, trichloro-
U209	1,1,2,2-Tetrachioroethane	U228	Trichloroethylene
U209	Ethane, 1,1,2,2-tetrachloro-	U234	1,3,5-Trinitrobenzene (R,T)
U2:0	Ethene, tetrachloro-	U234	Benzene, 1,3,5-trinitro-
U210	Tetrachloroethylene	U235	1-Propanol, 2,3-dibromo-, phosphate (3:1)
U211	Carbon tetrachloride	U235	Tris(2,3,-dibromopropyl) phosphate
U211	Methane, tetrachloro-	U236	2,7-Naphthalenedisulfonic acid,3,3'-[(3,3'-
U213	Furan, tetrahydro-(I)		dimethyl[1,1'-biphenyl]-4,4'-
Ų213	Tetrahydrofuran (I)		diyl)bis(azo)bis[5-amino-4-hydroxy]-, tetrasodium salt
U214	Acetic acid, thallium(1+) salt	U236	Trypan blue
U214 ·	• •	U237	2,4-(1H,3H)-Pyrimidinedione, 5-[bis(2-
U215	Carbonic acid, dithallium(1+) salt		chloroethyl)amino]-
U215	Thallium(I) carbonate	U237	Uracil mustard

#### SOURCE CODES

Code	Waste source	Code	Waste source
CTEAT	NING AND DEGREASING	A54	Oil changes
		A55	Filter/Battery replacement
A01	Stripping	A.56	Discontinue use of process equipment
A02	Acid cleaning	A57	Discarding off-spec material
A03	Caustic (Alkali) cleaning	A58	Discarding out-of-date products or
A04	Flush rinsing		chemicals
A05	Dip rinsing	A.59	Other production-derived one-time and
A06	Spray rinsing		intermittent processes
A07	Vapor degreasing	A60	Sludge removal
A08	Physical scraping and removal		
A09	Clean out process equipment	REMEDIATION DERIVED WASTE	
A19	Other cleaning and degreasing	A61	Superfund Remedial Action
		A62	Superfund Emergency Response
SURF	ACE PREPARATION AND FINISHING	A63	RCRA Corrective Action at solid waste
A21	Painting		management unit
A22 -	Electroplating	A64	RCRA closure of bazardous waste
A23	Electroless plating		management unit
A24	Phosphating	A6 <b>5</b>	Underground storage tank cleanup
A25	Heat treating	A69	Other remediation
A26	Pickling		
A27	Etching	POLL	UTION CONTROL OR WASTE
A29	Other surface coating/preparation		TREATMENT PROCESSES
	(Specify in Comments)	A71	Filtering/screening
			Metals recovery
PROC	ESSES OTHER THAN SURFACE	Ą73	Solvents recovery
	PREPARATION	A74	Incineration/Thermal treatment
A31	Product rinsing	A75	Wastewater treatment
A32	Product filtering	A76	Sludge dewatering
A33	Product distillation	A77	Stabilization
A34	Product solvent extraction	A78	Air pollution control devices
A35	By-product processing	A79	Leachate collection
A36	Spent catalyst removal	A89	Other pollution control or waste treatme:
A37	Spent process liquids removal		
A38	Tank sludge removal	OTHE	R PROCESSES
A39	Slag removal	A91	Clothing and personal protective
A40	Metal forming		equipment
A41	Plastics forming	A92	Routine cleanup wastes (e.g., floor
A49	Other processes other than surface		sweepings)
	preparation (Specify in Comments)	A93	Closure of management unit(s) or equipment other than by remediation specified in codes A61 - A69
PRO	DUCTION OR SERVICE DERIVED ONE-	494	•
	TIME AND INTERMITTENT PROCESSES	499	Laboratory wastes Other
ASI	Leak collection		

Cleanup of spill residues

A53

Code	Waste description	Code	Waste description	
	LAB FACKS	<b>B</b> 116	Leachate	
		B117	Waste liquid mercury	
LAB PACKS - Lab packs of mixed wastes, chemicals, lab wastes		B119	Other inorganic liquids (Specify in Comments)	
B001	Lab packs of old chemicals only	ORGA	NIC LIQUIDS - Waste that is primarily	
B002	Lab packs of debris only	organic and is highly fluid, with low inorganic sol content and low-to-moderate water content		
B003	Mixed lab packs			
B004	Lab packs containing acute hazardous	2000	C	
	wastes	B201	Concentrated solvent-water solution	
B009	Other lab packs (Specify in Comments)	B202	Halogenated (e.g., chlorinated) solvent	
		B203 B204	Non-halogenated solvent	
	LIQUIDS	B20 <del>4</del>	Halogenated/non-halogenated solvent mixture	
		B205	Oil-water emulsion or mixture	
	GANIC LIQUIDS - Waste that is primarily	B206	Waste oil	
inorganic and highly fluid (e.g., aqueous), with low suspended inorganic solids and low organic content		B207	Concentrated aqueous solution of other organics	
B101	Aqueous waste with low solvents	B208	Concentrated phenolics	
B102	Aqueous waste with low other toxic	B209	Organic paint, ink, lacquer, or varnish	
	organics	B210	Adhesives or epoxies	
B103	Spent acid with metals	B211	Paint thinner or petroleum distillates	
B104	Spent acid without metals	B212	Reactive or polymerizable organic liquid	
B105	Acidic aqueous waste	B219	Other organic liquids (Specify in	
B106	Caustic solution with metals but no cyanides		Comments)	
B107	Caustic solution with metals and cyanides		SOLIDS	
B108	Caustic solution with cyanides but no metals			
B109	Spent caustic		INORGANIC SOLIDS - Waste that is primarily morganic and solid, with low organic content and low-to-moderate water content; not pumpable	
B110	Caustic aqueous waste			
B111	Aqueous waste with reactive sulfides		-	
B112	Aqueous waste with other reactives (e.g.,	B301	Soil contaminated with organics	
	explosives)	B302	Soil contaminated with inorganics only	
B113	Other aqueous waste with high dissolved solids	8303	Ash, slag, or other residue from incineration of wastes	
B114	Other aqueous waste with low dissolved	9334	Other "dry" ash, slag, or thermal residue	
	soli <b>ds</b>	ä. :	Dry lime or metal hydroxide solids	

Code	Waste description	Code	Waste description	
B306	*Dry lime or metal hydroxide solids not *fixed*	B502	Lime sludge with metals/metal hydroxide sludge	
B307	Metal scale, filings, or scrap	B503	Wastewater treatment sludge with toxic	
B308	Empty or crushed metal drums or		organics	
	containers	B504	Other wastewater treatment sludge	
B309	Batteries or battery parts, casings, cores	B505	Untreated plating sludge without cyanides	
B310	Spent solid filters or adsorbents	B\$06	Untreated plating sludge with cyanides	
B311	Asbestos solids and debris	B507	Other sludge with cyanides	
B312	Metal-cyanide salts/chemicals	B508	Sludge with reactive sulfides	
B313	Reactive cyanide salts/chemicals	B509	Sludge with other reactives	
B314	Reactive sulfide salts/chemicals	B510	Degreasing sludge with metal scale or	
B315	Other reactive salts/chemicals		filings	
B316 B319	Other metal salts/chemicals	B511	Air pollution control device sludge (e.g., fly ash, wet scrubber sludge)	
<b>1</b> 313	Other waste inorganic solids (Specify in Comments)	B512	Sediment or lagoon dragout contaminated with organics	
ORGANIC SOLIDS - Waste that is primarily		B513	Sediment or lagoon dragout contaminated with inorganics only	
conten	organic and solid, with low-to-moderate inorganic content and water content; not pumpable		Drilling mud	
			Asbestos slurry or sludge	
B401	Halogenated pesticide solid	B516	Chloride or other brine sludge	
B402	Non-halogenated pesticide solid	B519	Other inorganic sludges (Specify in	
B403	Solid resins or polymerized organics		Comments)	
B404	Spent carbon			
B405	Reactive organic solid	ORGANIC SLUDGES - Waste that is prima organic with low-to-moderate inorganic so content and water content, and pumpable		
B406	Empty fiber or plastic containers			
B407	Other halogenated organic solids (Specify			
	in Comments)	B601	Still bottoms of halogenated (e.g.,	
B409	Other non-halogenated organic solids (Specify in Comments)		chlorinated) solvents or other organic liquids	
	SLUDGES	8602	Still bottoms of non-halogenated solvents or other organic liquids	
		B603	Oily sludge	
INORGANIC SLUDGES - Waste that is primarily		B604	Organic paint or ink sludge	
inorgan	nic, with moderate-to-high water content and	B605	Reactive or polymerizable organics	
low organic content, and pumpable		B606	Resins, tars, or tarry sludge	
B <i>5</i> 01	Lime sludge without metals	B607	Biological treatment sludge	

Code	Waste description	Code	Waste description					
B608	Sewage or other untreated biological sludge							
B609	Other organic sludges (Specify in Comments)							

#### GASES .

INORGANIC GASES - Waste that is primarily inorganic with a low organic content and is a gas at atmospheric pressure

B701 Inorganic gases

ORGANIC GASES - Waste that is primarily organic with low-to-moderate inorganic content and is a gas at atmospheric pressure

B801 Organic gases

DC		SIC		SIC	
004	Industry	Code	ndustry		neustry
	AGRICULTURE	1944	Subser area	2023 2024	Dry, condensed, evaluated products are cream and frazen deserts
		1081	Fernaday area, except variableM lights mining services	2024	Fluid Malt
GRIC	ULTURAL PRODUCTION—CROPS	1064	Uranium, radium, vanadium eres	2032	Canned specialists
2	Wheel Rice	1000	Metal ares, nes	2033	Cannos fruita ene vegetables
	Sam			2034	Denyerwed funts, regetation, south
•••	Sovoeers		MINING	2026 2027	Pictors, souces, and saled dressings Frazen fruits and vegetables
	Cash grains, nec	1221 1222	Stummeus cost and lights - suffess Stummeus cost - underground	2036	Frazen specialtes, nec
131	Cotton	1231	·	2041	Flour and other grain mill products
32	Tobacca Sugar cane and sugar beets	1241	Cast mining services	2043	Coreal breastest foods
• 34	nen parason			2044	Pice miling
138	Freid croos, except cash grains, nec		ND CAS EXTRACTION  Crude personum and natural gas	2045 2046	Prepared flour mixes and doughs Wet carn milling
161	regetables and melane	1311	Netural gas figures	2047	Deg and cat food
171	Serry cross Graess	1361		2046	Propared foods, nes
173	Tree num	1342	Oil and gas audioration contact	2061	Bread, case, and rested products
174	Citrus FURS	1386	Oil and gas field services, nes	2082	Coorde and oracions Fregor beauty products, except bree
175	Deciduous tree fruits		AETALLIC MINERALS, EXCEPT FUELS	2061	Rest cane sugar
179	Fruits and the rule. not	1413	Omenser sere	2042	Care sugar refining
181	Omemental numery products Food cross grown under cover	1422	Cruened and bressn limesters	5063	Seet sugar
191	General terms, primerly cross	1423	Cruence and broken grands	2064	Candy and other confectionery prod
-	-	1429	Cruenes and brown stens, nos	2066 2067	Chanada and cooks products Channe gum
S POS	CULTURAL PRODUCTION—LIVESTOCK	1448	Construction sand and grant industrial sand	2066	Safet and readed nuts and seeds
211	Good castle 'soulists	1456	Keein and had clay	2074	Cottonecod of mile
2'1 2'1	Seef cattle, encopt foodfolk Mags	1466	Clay and related manerals, nee	2075	Sayteen at mile
2214 2214	Shoop and goots	1474	Potest, sees and berste minerals	2076	Vegetains of mills, nec
218	General Addition, net		Photonical rect	2077	Arumas and marine has and oils
241	Davy ferms	1478	Chemical and fertilizer mining, not	2079	Egible tals and side, not Mad bevertages
2251	Broxer, fryer, and reaster chickens	1481	Nonmetalle minerale sentres Mineralengue nonmetalle minerale, nes	2003	Had
223	Chicago aggs	-		2064	Winds, brendy, and brandy source
2257	Turkky and furkky oggo Roughy nashanas			3065	Classics and blandes fever
238	Poutry and eggs, nee		CONSTRUCTION	2006	Bottod and served ask errors Fewering extrasts and syruce. ***
3271	Fur-bearing animals and rabbits		TRAL BUILDING CONTRACTORS	2067 2081	
2272	HOUSE BUT BEING CONTACT	1921		2000	Fresh or frecen propered fish
2273 2278	Animal aquacidhara Animal speciallies, nes	1522	- ·	3006	Roased cafee
327	General forms, primarily shares	1551		2007	Manufactured ine
	,	1541		2094	Meseron and stepholic Food proportions, not
	CULTURAL SERVICES	1542	Nanvesidential construction, not		
37:1		HEAT	Y CONSTRUCTION, DICLUONIQ SUILDINGS	TOE	LCCG PRODUCTS
יברנ בברנ	- · · · · •	1611	Highesty and street construction	_	Cigarettes .
3723	• • • • • • • • • • • • • • • • • • • •	1 (22)		2121	Cigars
3724		1623		2131	Chausing and amoung tobacco Tebases somming and rednying
2741		1 (66)	Heavy construction, not	2141	in the second second
3742		100	PLAL TRADE CONTRACTORS	TEXT	TLE MILL PRODUCTS
0751		1711		2211	Breakwaven faktis mills, cellen
0781		1721	Perrang and paper hanging	2221	Breadwayen faiths male, man-made
2762	Form management services		Electrical work	2231	Name to the make
	Condecage coursely and planning		Mesonny and other stenowers Plessonne, enymes, and insulation		Warnen's heavy, exists secus
	Caun and garden concess		Terrace, tile, markie, messis starts	2292	•
3783	Omenania shrub and the services	1751	_	2253	
FOR	ESTRY	179	Fleer leying and floer work, nee	2254	
3611	Timper tracks	1761		2257 2254	
	Forest products		Congress work Water well driving	2230	
185	Foresty services		Shrings at an execution	2361	
me:	HING, HUNTING, AND TRAPPING	-	Glass and grazing more	2200	
	2 Finden		Exemples were	2300	Freehing plants, not Carpote and rugs
391	3 Shedien		Wresting and demoision work		ARM SELLENG WAR
	Messeleneous manne products		Indexing building columns, not   Special triple contractors, not		Throwing and winding mile
	Figs Agriculture and Street Street	178	A CONTROL OF THE PROPERTY CONTROL OF CONTROL	2364	Throad mile
<b>JB</b> 7	1 moving, traceing, game presegation			2294	
			MANUFACTURING	2204	
	MINING			2297 2200	
	•		DO AND KINGRED PRODUCTS	2300	
	TAL MINING		1 Most passing plants 3 Sourcegos and other processed mosts		<del>-</del>
101	t risks area	٠٠.			

Note: nec = not elsewhere classified.

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Note: nec = not elsewhere classified.

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	umbing fisture fittings and from	35 <b>00</b> 35 <b>02</b>	Service industry mechinery, nee Careurotors, pistons, migs, venee	3843	Comes equipment and suggests X-ray appearatus and tubes
	eding equipment, except electric springs of the structure metals	3543	Fluid power cynnders and actuators	3845	Electromedical seviement
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	ofs, nuts, mers, and washers	3621 3624	Motors and generators Carbon and grashine preducts	3831	Museum merumana
	on and steel forgings lorriemous forgings	3625	Rolleys and industries controls	3842	- · -
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	roune and closures	3631 3632	Household cooling equipment Household refrigerators and freezers	3848	Searcing and ethicitic goods, not Pens and mechanical sensition
	letar stampings, net Tebng and powering	3633	Household leundry equipment	3002	Lead pencils and art goods
	letal coating and alled services	3634	Electric housewares and fans	3863	Mertang devices
	mos arms ammunison	3636	Household vacuum clashers	3045	Carbon separ and inted risbans Costume issuery
	mmunition, except for small arms, nec	3639	mousenoid appliances, nec Eastre lames		Femore, bullets, reedles, and pine
	imali arms Indinance and accessines, nec	3643	Current-carrying winning devices	3001	
-	ACTUAL VALVES	3644	Noncurrent-carrying mining devices	3005	Signs and advertising specialities
	fund power verves and hose fittings	3646	Rendertiel lighting fistures	3005	Surei custose finas empresas and
	ical sarrigs, except sare	3648 3647	Commercial lighting fishings Vehicular lighting equipment	-	Herd surface flaar eavenings, nes Menufacturing industries, nes
	laves and pipe fillings, net Vire service	3648	Lighting equipment, net		
	Meditionedus febricated wire products	3051	Household audio and video oculament	-	
167 M	Actor fool and vadi	3052	Prerecorded records and (appl)	11	ransportation and utiliti
	apricised pies and fittings	3661	Telegrame and telegraph apparatus  Redie and TV communication equipment	RAIL	DAD TRANSPORTATION
<b>459</b> F	denominal mates products, not	3000	Communications equipment, nes		Retroads, line-hour operating
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	during machinery	3678 3679	Electronic comectors Electronic comecnents, het	4131	
	Oil and gas field machinery Bevesors and moving diametry	3881	Signage befores	4142	
	Conveyors and conveying sourcement	3882	Primary betteries, dry and well	4151	
	Holets, cranes, and monorale	3004	Engine electrical equipment Magnetic and optical recording media.	4173	Bus terminal and service facilities
	ndustrial trucks and tractions	3665 3665	Engineer and opinion and supposes, nee	TRUC	DING AND WAREHOUSING
	Machine spose, metal sulting Types Machine spose, metal ferming Types		,		Local Studieng, without marage
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	Ar and gas compressive	3700 3700			Water transportation of freight, not
	Scores and fant Faceaging machinery	3706		4461	Door tot passanger trans., except farry
3544	Speed changers, areas, and goors	3790	Transportation equipment net	4482	- ··
	Industrial furnices and evens	-	NUMERTS AND RELATED PRODUCTS	4489	
			Search and navigation equipment		Towng and tugbed server
			LABORMONY SOCIETIES SIC TURNING	4463	
	- T-1		Emergramental controls	4400	Water transportation services, not
			Fraces control instruments Fluid maters and count no services		
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3561	Automatic vending machines		Ophical instruments and entire		
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} •	STID Industrial and partiang spread pages?	5718 Minamerican remarkamentary stores
4731 Fragre termentation arrangement	5131 Pleas goods and religion	5722 Hausenberg agentary agents
4741 Persel of religious care	5136 Man's and bayer stelling.	573d Company and sphengy stores
4736 Presection and fraud facilities	5130 Formula	5728 Named and provinced type starts
CONTRACTOR STANDS	5141 Granner, general fine 5142 Passagest Frager heads	
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AND COMMUNICATION SPRINGS, AND	STAR Greaters and respect products, rate STES Great and fines bears	
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4911 Chemita sarvanas		Shirt James Commercia
AND CAN PERSONAL AND	SIGN Character and affect greature, res	Shell ready, wy, and game areas
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SOOT Lamber, process, and military	STIT CHARTMAN BOWN	1
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SC Code	'noutry	Code	industry	SCC	Industry
	RITY AND COMMODITY BROKERS	7323	Credit recording services	HEAL	TH SERVICES  Offices and clinics of medical socions
5211	Security brokers and dealers Commodify contracts brokers, dealers	7331 7334	Direct mest asherbeing services Photocopying and duencabing services	8021	Offices and clinics of demists
5231 5231	Security and commonly exchanges	7336	Commercial shotography	8031	
212	megment stace	7336	Commercial art and greater dough	8041 8042	
5206	Security and commodify services, nec	7336 7342	Secretarial and sout resorting Distributions and sout control connects	8043	Office and clinics of podestries
MSUI	RANCE CARRIERS	7546	Building maintenance servess, nec	80-40	Offices of health gractitioners, nec
4311	ute insurance	7362	Medical countriest rental	808: 8082	
5321	Accident and health insurance mospital and medical service plans	7353 7360	Heavy construction couldment rental Equipment rental and leasing, not	8086	
5324 5331		7381	Emeloyment agencies		General modical and surgical hosoitals
6361	•	7363	Hole supply services	8083 8088	Psychiatric hospitate Sauciaty nospitate, except psychiatric
6361	Title insurance	7371 7372	Corrector programming services Programmed software	8071	
5371	Pension, neath, and siefere funds reurence carriers, nes	7373		8072	<del>-</del>
		7374	Data processing services	NORTH THE	
	RANCE AGENTS. BROKERS, AND SERVICE	7375		ACRES	Kidney disysts corners Souccety outpetent clinics, nec
5411	nsurance agents, brossis, and service	7376 7377	Computer rental and leasing		Health and allied services, nes
MEAL	STATE	7378	Computer maintenance and report		
	Honraudentes building operators		Computer revised services, net		SERVICES
95'3		7381 7382		•111	Legal services
6514		7363	Name synerces	EDUC	ATIONAL SERVICES
8617		7384	Photofinaning laborationes	4211	Elementary and secondary schools
8619	Real property lessors, net	7300	Business services, nec	8221 8221	Calegos and universities  Juner eatleges
<b>66</b> 41	Real estate opens and managers This assirant offices	AUTO	MOTIVE REPAIR, SERVICES, AND PARKING		Librarios
8562	Subdividers and developers, net		Truck rantal and leasing, no others	1243	Cass processing schools
***	Commeny supplieders and developers	7514		524	Supress and secretarial schools Venational schools, not
		7515 7518	Pessonijer car teating Utility trader rental	1244 1288	Sphools and educations services, rec
	DING AND OTHER INVESTMENT OFFICES  Bark holding companies	7318	Adametric pertang		
4718	<del></del>	7538	Top and body repair and paint shops		AL SERVICES
	Management investment, open-and	7535	Auto exhaust system repair sheet	8322 8331	Individual and family services Jap training and related services
6725		7534 7536	The retreating and reser those Augumphie glass resecoment those	8381	Child day care services
6733	Educational religious, etc. truets Truets, nec	7937	Automobile transmission repor sheet	8361	
6792	Oil royally traders	7536	General automobile report throat	1300	Secret services, nes
	Powers owners and leasons	7538) 7542	Automobie reper shops, nes Car weshes	MUSE	TUMB, BOTANICAL, ZOOLOGICAL GARDENS
6796		75-40	Automotive services, red		Museums and art gallenes
•				1422	Botanical and sollogical gardens
	SERVICES		ELLAMEOUS REPAIR SERVICES  Reads and total minute.	MON	MERCHAP ORGANIZATIONS
	SERVICES	7923	Refrigeration service and most		Business associations
HOT	TLE AND OTHER LOOGING PLACES	7000	Bearing repair shees, net		Professional organizations Labor organizations
7011		7631 7641	Weigh, discit, and jovery repair. Reveniessery and furniture repair.	20031 20041	
7021		7000	Weiding repor		Political organizations
*333		7004	Armeture reserveing shops	8881	
7041	Membershie-basis organization holets	7000	Regar services, net		Membershie organizations, nes
-	SONAL SERVICES	MOT	ION PICTURES		MEETING AND MANAGEMENT SERVICES
	Found laundred, family and commercial		Metten proture and reses production		Engineering services Arendeeburgs services
	Garment pressing and desirers' agents		Spreads affect to meltion pictures Median grature and cape drambunen		Support services
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	Local Contract Straight September 1998		DEMENT AND RECREATION BERVICES	8741	
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	1 Gener sheet		Theoretai producers and services Entertainers and emenantment groups	5744	
	1 Shop repair and shapshine shaps 1 Fungral service and cramesones	7933	Bauting centers	8746	Business consulting, nes
٠ <del>۵</del>			Sparts study, managers, and promoters	-	ATE HOUSEHOLDS
729	8 Meceterous personal services, nes		Resing, visitating track operation Physical Researchmen		Private hausthards
<b>SL</b> M	SHESS SERVICES		Public golf courses		
731	1 Assuming agencies	7962	Con-equipment amusement series		ACES, MBC Servens, nes
	2 Outdoor advertising services	7004	Amazoneri peris	-	

#### SIC CODES

(Continued)

SC Code Industry

SC Code Industry

SIC Code Industry

#### PUBLIC ADMINISTRATION

#### DIRCUTIVE LEGISLATIVE AND GENERAL

9111 Execute offices
9121 Legentate boses
9131 Execute and legentate communed
9130 General government, nec

#### JUSTICE, PUBLIC ORDER, AND SAFETY

9211 Courts 9221 Passes pro

9222 Legis courses and presentation 9225 Correctional methyleury

9224

\$220 Funds order and salety, nos

### PINANCE, TAXATION, AND MONETARY POLICY 9311 Prenes, lession, and mensery pasty

#### ADMINISTRATION OF HUMAN RESOURCES

\$411 Administration of educational programs \$421 Administration of public health programs

9441 Administration of social and management

programs \$461 Administration of votorgraf affairs

#### ENVIRONMENTAL QUALITY, AND HOUSING

9611 Air, water, and solid waste manager 9612 Land, mineral, whethe agreements

9651 Housing programs 9632 Urban and assimulate of

#### ADMINISTRATION OF ECONOMIC PROGRAMS

Agents THATTICK OF SCORCING PROTORNA 3611 Advan, of general connection programs 1921 Agentsian, advant, of temperation 1931 Regulation, advantabation of utilities 1941 Regulation of agricultural managing 1951 Space receives and technology

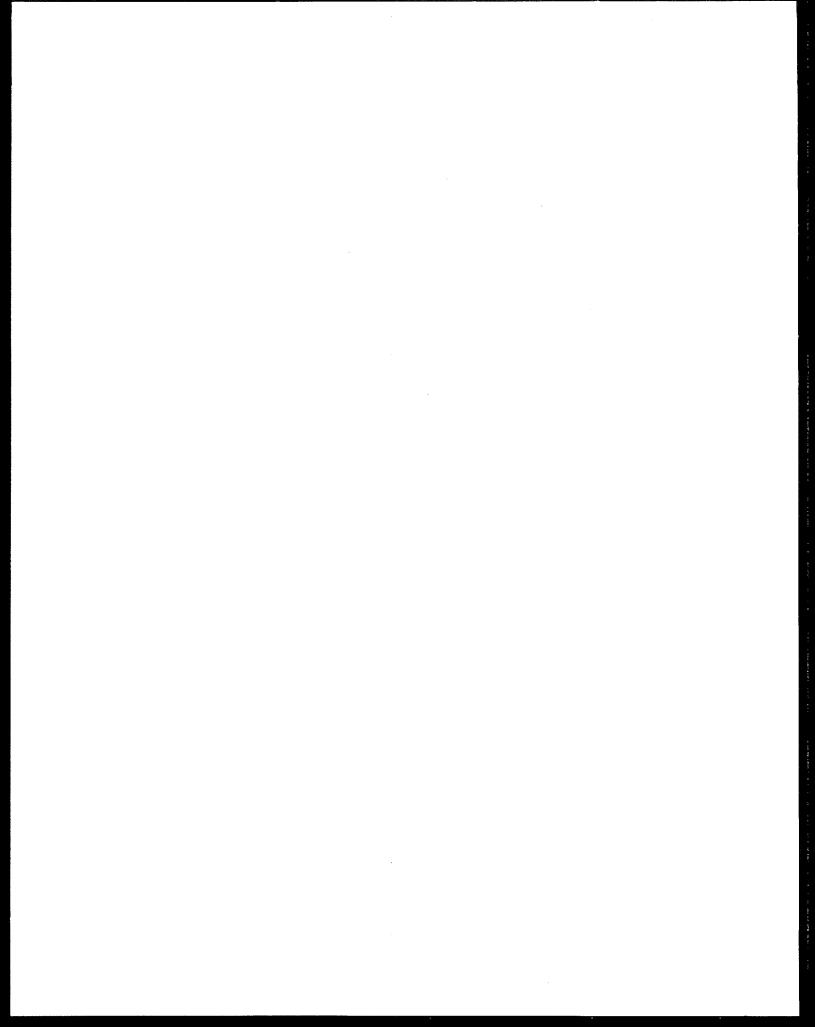
#### NATIONAL SECURITY AND SITEMATICINAL APPAINS 9711 Helland security 9721 International allows

#### **NONCLASSIFIABLE ESTABLISHMENTS**

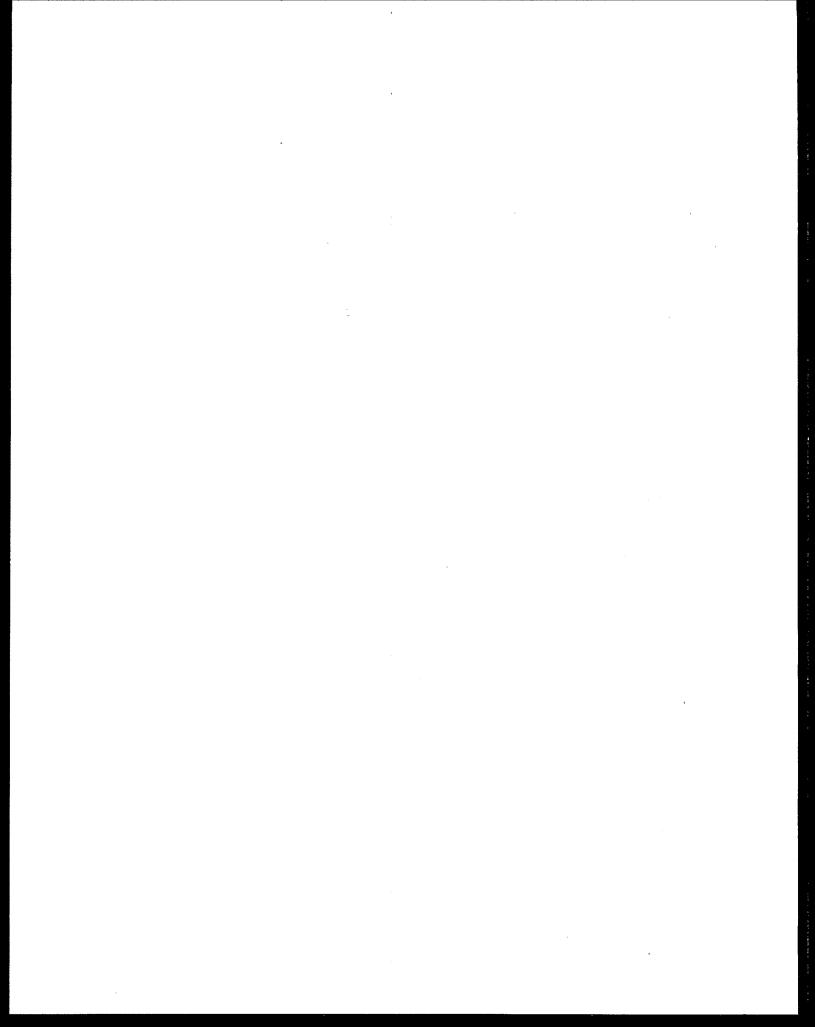
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#### **ORIGIN CODES**

Code	Origin Description
1	The hazardous waste stream was generated on site from a production process or service activity (including off-specification or spent chemicals).
2	The hazardous waste stream was the result of a spill cleanup, equipment decommissioning, or other remedial cleanup activity.
3	The hazardous waste stream was derived from the management of a non-hazardous waste stream.
4	The hazardous waste stream was received from off site and was not recycled or treated on site.
5	The hazardous waste stream was a residual from the on site treatment, disposal, or recycling of previously existing hazardous waste stream.



## APPENDIX 3 MATCHING PROCESS - WR/GM



This Appendix summarizes EPA's manual and automated merging of GM and WR Forms data for source characterization of the wastes managed off site. As described in Chapter 2, EPA is using WR Form data to characterize the management of the wastes combusted off site, and source data (i.e., SIC code of the process generating the waste, origin code, and source code) are not listed on this form. Source data are needed to determine how wastes were generated and identify waste minimization opportunities, and thus, it is important to link the WR forms to GM forms as accurately as possible in order to promote waste minimization for wastes managed off site.

EPA used both automated and manual techniques to merge WR data with corresponding GM data. The automated mapping was done for both combusted and fuel-blended wastes. The manual merging was done only for the combusted wastes. The first step in the merging process was the aggregation of data by routing information (i.e., generator and management facility IDs), RCRA waste codes, management system type, and waste form codes; these information elements are common to both GM and WR Forms. The merging was done in three stages:

- 1. Initial automated "exact" merging;
- 2. Manual matching of data for the top 50 wastestreams; and
- 3. Final automated "relaxed" merging.

These three steps are described below:

Step 1. Automated "Exact" Merging: The objective of the initial automated "exact" merge was to obtain waste stream-specific source information from the data without making any concessions in the matching criteria. This process resulted in a poor match of the common information between GM and WR Forms. Only 9,696 records, contributing a volume of 82.719 tons of combusted and fuel-blended wastes (including 1,104 records of combusted wastes with 65.200 tons, or about two percent of off-site combusted waste quantity), were matched in this stage.

Step 2. Manual Matching: Of the remaining non-match combustion records (25.257 records representing 1.1 million tons of combusted waste), the top 50 records accounted for about 40 percent of the total waste quantity. EPA used a manual process, as described in Attachment 1, to match the GM and WR forms' information. This process resulted in identifying source data for an additional 33 records contributing a quantity of 220,102 tons, or about 25 percent of the off-site combusted waste quantity. Moreover, this step provided some insight on how to set up decision rules for the final step.

Step 3. Automated "Relaxed" Merging: EPA aggregated the remaining non-matched records (107,552 records contributing 3.9 million tons, or 91 percent, of combusted and fuel-blended wastes) by routing information, RCRA waste codes, management system type, and waste

<sup>&</sup>lt;sup>1</sup> Fuel-blended wastes were later dropped from the analysis in order to focus on combusted wastes.

form grouping code (see Exhibit 1 of Attachment 1 for a list of the waste form groups). These records were electronically mapped to the GM forms data using routing (i.e., generator sending the waste and managing facility receiving waste are identical on the GM and WR forms) and RCRA waste code information. Where there were perfect matches for routing and RCRA waste code data, the source data were obtained from the GM Form. This process resulted in mapping of 14,891 records contributing 593,699 tons of waste.

EPA aggregated the remaining non-matched records (88,375 records) to make a smaller file for downloading into the PC environment. EPA's strategy for downloading was as follows:

- For records with quantities greater than or equal to 1 ton, keep all data as reported. There are 36,744 records representing 3,367,719 tons in this category.
- For records with volumes less than or equal to 1 ton, aggregate by receiving facility ID, generator state (i.e., first two characters of the generator ID), waste code(s), management system type, and waste form group code (see Exhibit 1). This results in losing information on individual generator ID and individual form codes. In this step, 51,631 records collapsed to 22,957 records representing 18,089 tons.

Overall, for combusted and fuel-blended wastes, the aggregation and merging results are as follows:

Waste Type	With Sou specific Ir	rce- iformation	Without S specific In	Source- oformation	Total		
	Records	Tons	Records	Tons	Records	Tons	
Combusted Wastes	7,716	474,363 (40%)	21,802	716,568 (60%)	29,518	1,190,931	
Combusted and Fuel- blended Wastes	24,587	896,520 (21%)	59,701	3,385,808 (79%)	84,288	4,282,328	

Please note that due to non-reporting of data in the GM Forms, even for perfectly matched records some of the source data could be missing. Thus, even after matching there are only 462.203 tons of waste with SIC code information and 437,394 tons of waste with source code information for off site combusted wastes.

#### Attachment 1. Manual Matching of GM and WR Forms for Selected Combusted Wastes

This attachment details the methodology and results of the manual matching done for selected combusted waste streams identified in the WR Forms that could not be electronically matched to the GM Forms. The primary objective of the matching process is to obtain source data (i.e., source, origin and SIC codes) for routinely-generated wastes identified in the WR Forms. The data obtained from the matching would be used in identifying and prioritizing the sources of combusted hazardous wastes.

There are about 30,000 WR Form records for which source data could not be obtained by automated "exact" matching and merging. The top 50 records from the non-matched WR Form records contribute to about 42 percent of the non-matched waste quantity and the top 1,300 records contribute to about 76 percent of the total non-matched waste quantity. Hence, the manual matching was done for the top 50 waste streams in order to address a significant percentage of the unmatched quantity and to develop and implement an automated programming logic for establishing source data for most of the remaining non-match records.

The first step of the manual matching was to ascertain that the waste routing information is accurate (i.e., EPA IDs reported in GM and WR Forms are identical). If this condition could not be fulfilled, it was assumed that the waste stream in the WR Form could not be matched to the GM Form. After the routing information was verified, then the data in the WR and GM Forms were assumed to be matched if they met at least two of the following three conditions:

- (1) The waste form codes of the waste stream in the GM and WR Forms indicate similar wastes. The groups of waste form codes that were considered to be similar for this analysis are given in Exhibit 1.
- At least one of the RCRA hazardous waste codes of the waste stream as reported in the GM Form is also reported in the WR Form.
- (3) The quantity of the waste stream reported in the GM Form is within 25 percent of the quantity reported in the WR Form.

Exhibit 2 presents the results of the manual matching. The analysis indicate that the matches between the WR and GM Forms were often poor. After using the broad assumptions outlined above, 33 of the 50 top non-matched WR Form wastes were matched. For these 33 waste streams, the summary findings are presented below:

- For 13 of the 33 waste streams:
  - the waste form codes were similar (based on Exhibit 1);
  - at least one of the EPA hazardous waste codes matched; and
  - the reported quantities were within 25 percent between the WR and GM Form data.
- For 11 of the remaining 20 waste streams:
  - at least one of the EPA hazardous waste codes matched; and
  - the reported quantities were within 25 percent between the WR and GM Form data:

- For six of the remaining 9 waste streams:
  - the waste form codes were similar (based on Exhibit 1); and
  - at least one of the EPA hazardous waste codes matched between the WR and GM Form data.
- For the remaining three waste streams:
  - the waste form codes were similar (based on Exhibit 1); and
  - the reported quantities were within 25 percent between the WR and GM Form data.

For the 17 WR Form wastes that could not be matched, listed are some of the reasons why the matching could not be done:

- For four waste streams, the WR Form identified a generator ID that is not in the BRS;
- For one waste stream, the WR Form identified a generator that had no GM Forms;
- For seven waste streams, the GM Forms of the generators identified in the WR Forms did not report snipping wastes to the receiver in 1991; and
- For five waste streams, though the GM Forms did report shipping some wastes to the receiver, none of the waste streams could be matched since, for each of these waste streams at least two of the three conditions outlined above were not fulfilled between the WR and GM Forms.

Barring a few exact matches, one of the three common parameters (i.e., the RCRA hazardous waste codes, the waste form codes, or the quantities) are different between the WR and GM Forms for almost all of the manually matched waste streams. Even though the generator may have a better knowledge of their waste streams, for the final analysis of combusted wastes the information provided in the WR Forms on the three parameters will be used. This is because the receiver would be expected to determine the exact properties (and quantity) of the waste prior to treating the waste for several reasons including:

- to determine that right treatment system;
- to determine the exact price for treating the waste stream; and
- to avoid liability resulting from not treating the waste to LDR standards.

Note that several of the waste streams that were matched will be deleted for the final evaluation of combusted wastes. This is because the matching determined that these waste streams are not routinely-generated wastes (i.e., source codes were A61-A69 or origin codes were 2 or 5).

Exhibit 1: Groups of Form Codes That are Similar

The groups that were considered to be similar for the manual matching of the metal-bearing combusted wastes are listed below:

1) B20x and B219	2) B201 and B202
3) B201 and B203	4) B202 and B204
5) B203 and B204	6) B205 and B603
7) B301 and B401	8) B40x and B409

Similarly, the groups of similar form codes can be expanded to include:

9) B00x and B009	10) B1xx and B119
11) B103 and B105	12) B104 and B105
13) B106 and B110	14) B107 and B110
15) B108 and B110	16) B111 and B112
17) B205 and B206	18) B209 and B604
18) B206 and B603	19) B3xx and B319
20) B5xx and B519	21) B60x and B609

For a more detailed explantion of codes, see Appendix 2.

Exhibit 2: Top 50 Waste Streams in the WR Forms That Were Manually Matched to the GM Forms

085	Receiver's 1D (WR/GH)	Generator's ID (WR/GM)	EPA Haz Codes (WR)	EPA Haz. Codes (GH)	Form Code (WR)	Form Code (GM)	SIC Code (GH)	Source Code (GM)	Origin Code (GM)	Quantity (WR) (short tons)	Quantity (GH) (short tons)
1	1XD083472266	TXD058265067	D001 D018 D035 f003	D001 D002	-	8207	2869	A35	ı	36,916	45,800
2*	KSD031203318	M109806152 <b>98</b>	DOO1 DOO5 DOO6 DOO7 DOO8 DO18 DO26 DO35 FOO1 FOO2		B219					, 148	43,600
3*	ARD981512270	MID980615298	D001 D004 D005 D006 D007		B20 <b>4</b>					21,086	
4	PAD002389559	HJD002454544	F003 F005	D001 D035 D038 F003 F005	B219	B203	7389	A49	5	20,632	23,137
5.	VAD042755082	VAD098443443	D001 D018 D043 F001 F002 F003 F004 F005		B204					15,510	
6	MOD050232560	IND980590947	D001 F002 F003 F005	D001 F001 F003 F005	B202	8204	2869	A89	5	11.400	
<i>i</i> •	M0D054018288	₩1D000808824	D001 F001 F002 F005	·	B204			703		11,420	132
8	ARD981512270	TXD000/42304	D001 D005 D006 D007 D008	f002-9 f019 K019 K030 K052 several U codes	B204	B203	7389	<b>A</b> 89	1	10,929 10,883	10,892
9.	PADOO2389559	NJD002182897	F003 F005	•	B219			<del></del>		10.016	
10	ARD981512270	LAD079464095	D001 D004 D005 D006 D007	D001 D004 D005 D006 D007	B204	B219	8999	A99	5	10,016 9,003	9,061
	IND005081542	OHD005051826	D001 D007 D008 D018	D001 K048 K049 K051	B204	B205	2911	A89	1	9.555	
12	MGD054018288	1ND093219012	D001 D004 D007 D009 F001 F002 F003 F004 F005 K048	F001 F002 F003 F004 F005	B204	B204	9999	<b>A9</b> 9	5	7,416	8,494 7,113
13*	ARD981512270	ARD981057870	D001 D005 D006 D007 D008		B407					7,323	
14	KSD031203318	IND000646943	D001 D007 D008 D018 D022 D026 D027 D028 D033 D036	0001 D002 D005-8 D018 D019 D022 F001-5 K052	B219	B204	9999	A99	4	7,002	6,641

Exhibit 2: Top 50 Waste Streams in the WR Forms That Were Manually Matched to the GM Forms (continued)

OBS	Receiver's ID (WR/GM)	Generator's ID (WR/GM)	EPA Haz. Codes (VR)	EPA Haz Codes (GM)	Form Code (WR)	Form Code (GM)	SIC Code (GM)	Source Code (GM)	Origin Code (GM)	Quantity (WR) (short tons)	Quantity (GM) (short tons)
15	PRD980526115	PRD981182421	F001 F002 F003 F005		B204					6,953	
	LAD008161234	TXD008090409	D001 D028 F037 F038	K048 K049	8205	B603	2911	A89	11	6,785	7,164
17	1ND072040348	1ND005460209	D001 F003 F005	F008	B201	B203	2819	A49	5	6,774	8,065
18	MOD050232560	MOD029729688	D001 D004 D005 D006 D007 D008 D009 D010 D011 D016		B202					6,591	
19	FLD000737312	GAD093380814	F001 F002 F003 F005	D001 D004 D006 D008 D011 F001-4 U's	B204	B204	2869	A89	5	6,442	640
20	ARD9815122/0	ARD981057870	0001 0005 <b>0006 0007</b>	D001 D005-9 D026 D035 D040 F001-5 K027	B204	B219	9511	A99	5	6,346	5,230
21.	MUDJ54018288	1ND984866541	D001 F003 F005		B204					6,065	
55.	M00050232560	W10000808824	D001 F001 F002 F003 F005.		B202					5,126	
23.	FL0004059085	FLD000737312	0001 F001 F002 F003		B204				<u> </u>	5,074	
24	ARD981512270	TXD046844700	D001 D004 D005 D006 D007	0004-20 F001-3 F005	B204	B204	7389	A73	5	4,936	5,180
25	MOD050232560	TMD991279480	D001 D007 D008 F001 F002 F003 F005	D001	B202	_	9999	•	1	4,866	4,904
26	M0D050232560	OKD058078775	K048 K049 K051	K048 K049 K051	B202	B504	2911	A75	5	4,794	5,038
27	KSD031203318	TXD000742304	D001 D004 D005 D006 D007 D008 D009 D010 D016 F001	D002 D004-8 D010 F001-8 F019 K,P,U codes	8219	B207	7389	A71	1	4,743	194

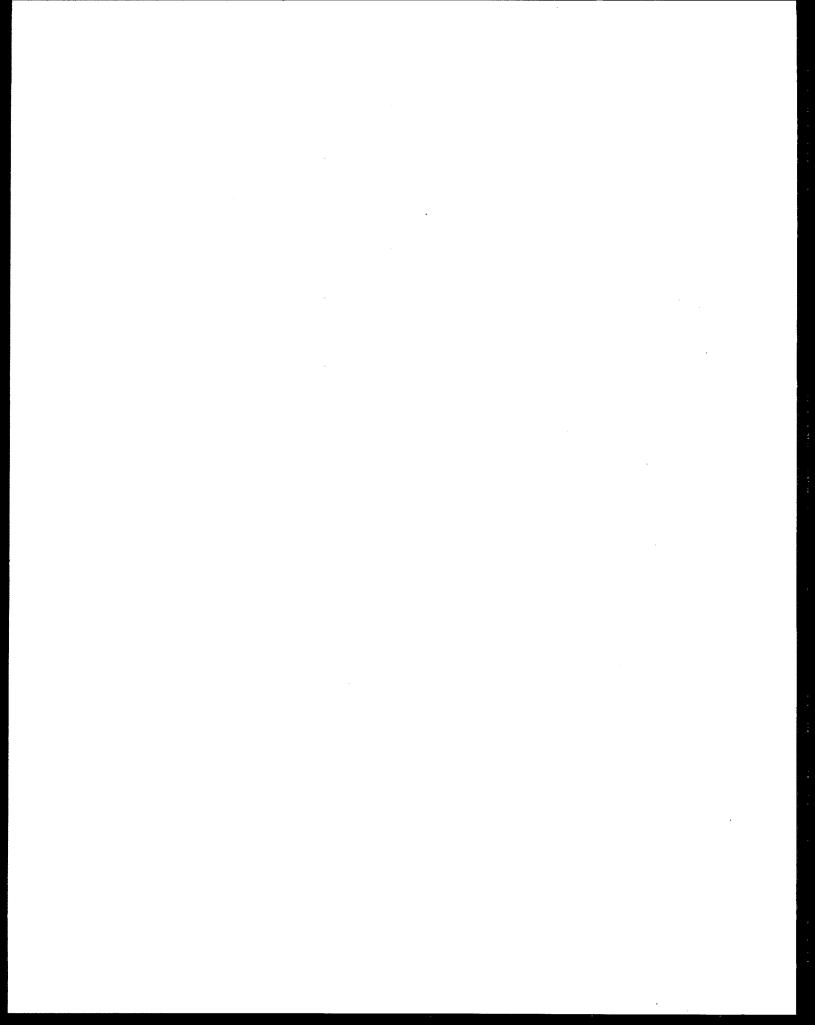
Exhibit 2: Top 50 Waste Streams in the WR Forms That Were Manually Matched to the GM Forms (continued)

100	A CONTRACT OF THE PARTY OF THE										
085	Receiver's ID (WR/GH)	Generator's ID (WR/GM)	EPA Haz. Codes (WR)	EPA Haz. Codes (GH)	Form Code (WR)	Form Code (GH)	SIC Code (GM)	Source Code (GH)	Origin Code (GH)	Quantity (WR) (short tons)	Quantity (GH)
28	LAD981057706	ALD070513767	D001 D004 D005 D006 D007 D008 D010 D011 D018 D035	D001 F001-3 F005	B204	8204	2899	A89	1	4,531	(short tons) 6,961
29	1ND005081542	1ND980590947	D001 F001 F002 F003	D001 F001-3 F005	B204	B204	2869	<b>A</b> 89	5		_
30	IND005081542	OHD004274031	D005 D006 D008 F001	D001 D005-8 D018 D035 D039 F001-3 F005	8204		4953	-	-	4,427	5,180 4,360
31	SCD036275626	NCD991278276	D001		B403						
32	MOD050232560	ILD066918327	D001 D018 D035 F001 F002 F003 F005	D001 F001-3 F005	B202	B201	7389	A19	5	4,320 4,313	5,318
33	H0D054018288	11 0087157251	D001 F002 F003 F005	D001 F001-3 F005	B204	B204	4953	A89			
34	1.A0000161234	LAD000812818	D001 D035	D001 D035	8203	B219	2869	A35	5	4,297	4,245
35°	\$XD006451090	LAD058530510	D008				2003			4,001	3,866
36	IND005081542	OHD093945293	D001 D005 D006 D007		B204					3,844	i
37	FXD981153711	TXD051161990	K051	K051		B603	2011			3,775	
38	IND005081542	OHD005046511	D001 D018 K048 K049	D001 K048 K049 K051	B204	- 5003	2911	A89		3,687	4,098
39	LAD981057706	TXD000742304	D001 D005 D006 D007	D002 D004-8 D010	B206		2911	-		3,669	3,477
		· · · · · · · · · · · · · · · · · · ·	D008 F001 F002 F003 F004 F005	F001-8 F019 K,P,U	D200	B207	7389	A71	1	3,518	2,611
40	AL 0981020894	VAD065408692	D001 F003 F005	0001 0035 F003 F005	B403	B409	2052				
41	OHD004304689	W10059972935	D001 D005 D006 D007	F003		B602	3053 2821	A56 A73		3,465	14
	<del></del>	·	D008 F003 F005				5051	۸/3	3	3,410	3,566

Exhibit 2: Top 50 Waste Streams in the WR Forms That Were Manually Matched to the GM Forms (continued)

085	Receiver's	Generator's	EPA Haz. Codes (WR)	EPA Haz. Codes (GM)	Form Code (WR)	Form Code (GM)	STC Code (GM)	Source Code (GM)	Origin Code (GM)	Quantity (WR) (short tons)	Quantity (GM) (short tons)
42	. ID (WR/GM) M0D050232560	1D (WR/GM) 1LD005476882	K049	K049	8202	8603	2911	A75	3	3,316	3,469
43	IND005081542	OHD980700942	D001 D002 D005 D006	D001 D006-8 D018 D035 D039 D040 F001-3 F005	B204	B204	4953	*		3,296	- 2,311
44	KSD031203318	LAD000618256	D001 D005 D020 P004 P010 P018 P030 P041 P048 P050	D007 P001 P004-5 P010 P018 U,F,K codes	B407	8609	4953	A64	1	3,165	3,344
45	PAD002389559	PAD064375470	F003 F005	D001 D004-8 D010-11 F001-3 F005 organic D codes	8219	6204	4226	A89	5	3,127	2,122
46*	OHD981195779	OHD093945293	D001 D005 D006 D007 D008 D011 D022 D035 D039 F001							3,124	
47	VAD077942266	PAD064375470	D001 D004 D005 D006 D007 D010 D011 D018 D019 D021	D001 D004-8 D010-11 F001-3 F005 organic D codes	B204	B204	4226	A89	5	3,107	3,029
48	IND005081542	OHD980681571	D001 F001 F003 F005		B204			,		3,057	
49"	KSD031203318	ARD981057870	D001 D005 D006 D007 D008 D009 D019 F001 F002 F003		B407	·				2,999	
50	1LD098642424	NJD986578953	F001 F002 U061	F001 F002 U036 U061 U129	B301	8401	4225	A66	2	2,900	1,869

Waste streams that could not be matched.



### APPENDIX 4 DOCUMENTS USED AS SOURCES OF CONCENTRATION DATA



### APPENDIX 4: SOURCES OF INFORMATION ON CONCENTRATIONS OF LISTED WASTES

- "Regulatory Impact Analysis of Restrictions on Land Disposal of Certain Solvent wastes," Office of Solid Waste, U.S. EPA, November 1986. Exhibit A-15 was used: this exhibit provides summarized information for all F001-F005 wastes that were incinerated in 1983.
- "Regulatory Impact Analysis of the Land Disposal Restrictions for Third Third Scheduled Wastes, Final Rule," Office of Solid Waste, U.S. EPA, April 25, 1990
- Listing Background Document for Petroleum Refining Wastes, August 15, 1988. This source only provides summarized information for the industry and does not provide waste-stream specific data; it was used for identifying constituents.
- Best Demonstrated Available Technology (BDAT) Background Documents.
  These included:

BDAT Background Document for F001-F005 wastes (Final), June 15, 1989

BDAT Background Document for F024 wastes (Final), May 15, 1989

BDAT Background Document for F037/F038 wastes (Final), June 30, 1992

BDAT Background Document for K002 wastes (Final), May 8, 1990

BDAT Background Document for K013 wastes (Final), June 15, 1989

BDAT Background Document for K016, K018, K019, and K020 wastes.

(Final), August 15, 1988

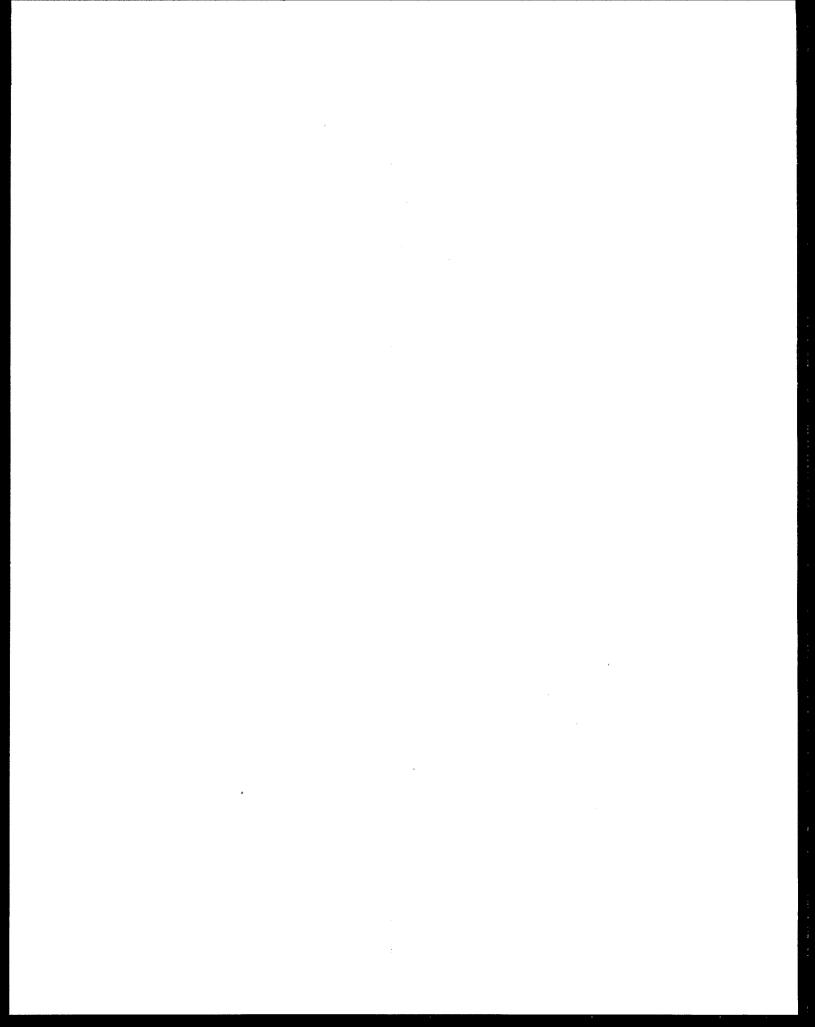
BDAT Background Document for K022 wastes (Final), August 15, 1990

BDAT Background Document for K028 wastes (Final), May 25, 1989

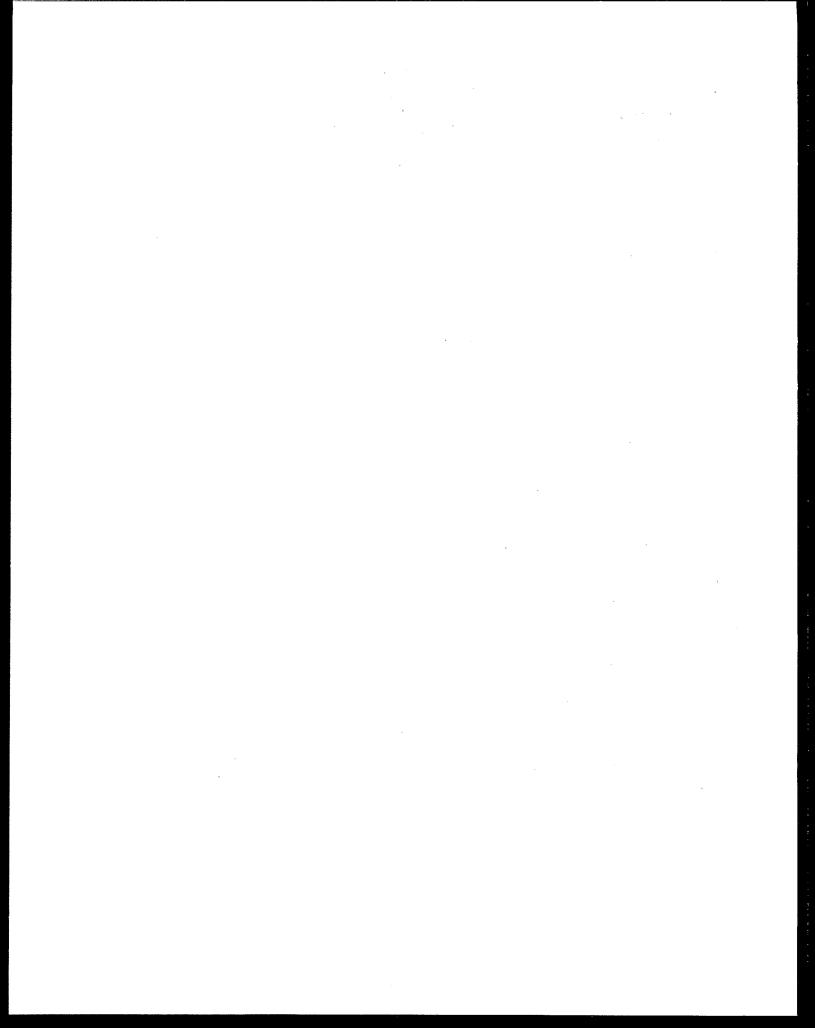
BDAT Background Document for K083 wastes (Final), May 8, 1990

BDAT Background Document for P063 wastes (Final), June 15, 1989

- \*Estimates of Waste Generation by the Organic Chemical Industry," Final Draft Report, U.S. EPA, December 7, 1987. This source was developed for the TC LDR RIA. It provides summarized information for the industry and does not provide waste-stream specific data.
- "Regulatory Impact Analysis: Proposed Standards for the Management of Used Oil," Office of Solid Waste, U.S. EPA. July 1985.



## APPENDIX 5 RETRIEVAL FROM GENSUR



#### APPENDIX 5: RETRIEVAL FROM GENSUR

For each waste stream combination, the four key attributes (i.e., RCRA code, SIC code, source code, and form code) were used to match the BRS information to corresponding waste streams in the Generator Survey (GENSUR). The constituent and concentration information for these matched GENSUR waste streams were assumed to be appropriate for the BRS waste stream combination. The following presentation details the procedure and assumptions used for matching the information.

First, an attempt was made to exactly match all four elements of the unique BRS combination with corresponding GENSUR combinations. Exhibit 1 provides the correlation for matching the form codes in the BRS to the waste description codes in the GENSUR. Exhibit 2 provides the correlation for matching the source codes in the BRS to the source codes in the GENSUR.

For those combinations for which an exact match could <u>not</u> be found in the GENSUR, the following assumptions were used (in descending order) for matching the four elements. If the usage of an assumption resulted in more than one GENSUR match for a particular BRS waste stream combination then the constituent information was averaged across all the matches for that combination.

- The RCRA codes exactly match and the BRS form code exactly matches the GENSUR waste description.
- 2) The RCRA codes exactly match and one or more of the other three elements exactly match.
- 3) 50 percent or more of the RCRA codes match and the form code exactly matches.
- 50 percent or more of the RCRA codes match and one or more of the other elements exactly match.
- 5) The RCRA codes exactly match.
- 6) One or more of the RCRA codes match and the form code exactly matches.
- 7) Any three of the matching elements exactly match.
- 8) Any two of the matching elements exactly match.

The median of the constituent concentration range given in GENSUR was used as the concentration for a given constituent (e.g., a concentration range of 6 in the GENSUR would translate to an estimated concentration of 50,000 ppm). These medians were used for averaging the concentrations of a given constituent across the waste streams in which the constituent occurs. Concentrations were weighted by quantities of <u>all</u> waste streams (i.e., final concentration was estimated by assuming that all the matched waste streams are mixed together).

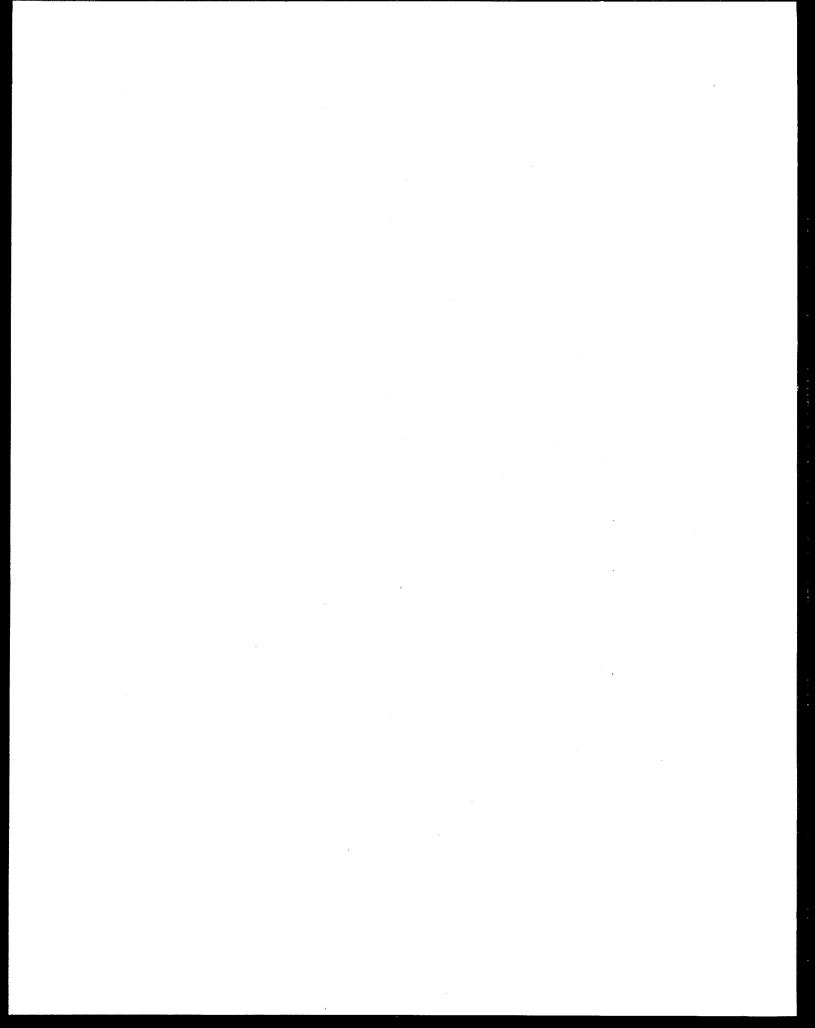
Exhibit 1: Form Codes for Top 200 Unique BRS Combinations

BRS Form Codes	Generator Survey - Waste Description Codes						
B101	B01						
B102	B02						
B105	B05						
B110	B10						
B111	B11						
B114	.B14						
B201	B58						
B202	B59						
B203	B60						
B204	B61						
B205	B62						
B206	B63						
B207	B64						
B208	B65						
B211	B68						
B212	B69						
B219	B70						
B301	B36						
B401	B80						
B403	B82						
B405	B84						
B407	B89						
B409	B90						
B503	B21						
B601	B71						
B602	B72						
B603	B73						
B606	B76						
B494, B597, B, blanks	B99						

Exhibit 2: Source Codes for Top 200 Unique BRS Combinations

BRS Source Codes	Generator Survey - Source Codes						
A09	S60						
A19	S03, S07, S12						
A31	S44						
A32	<b>S4</b> 0						
A33	S28						
A34	S27 .						
A35	S20						
A36 <sup>1</sup>	S15, S16, S26, S31-38, S48						
A37 <sup>1</sup>	S13-S16, S23, S25, S26, S30-S39, S41, S42, S45, S48						
A49	S13-S16, S23, S25, S26, S29-S39, S41, S42, S45, S48						
A54 <sup>1</sup>	S78						
A56	S61						
A57	S49						
A60 <sup>1</sup>	S16, S25, S41, S45, S48						
A71	S67						
A73	S74						
A74	S72						
A75	S73						
A89	S66, S69, S70, S71, S77						
A92	S60, S62, S78						
A99	S78						

<sup>&</sup>lt;sup>1</sup> There is a greazer degree of uncertainty in these code translations. Please use these translations with a lesser degree of confidence as compared to the other code translations.

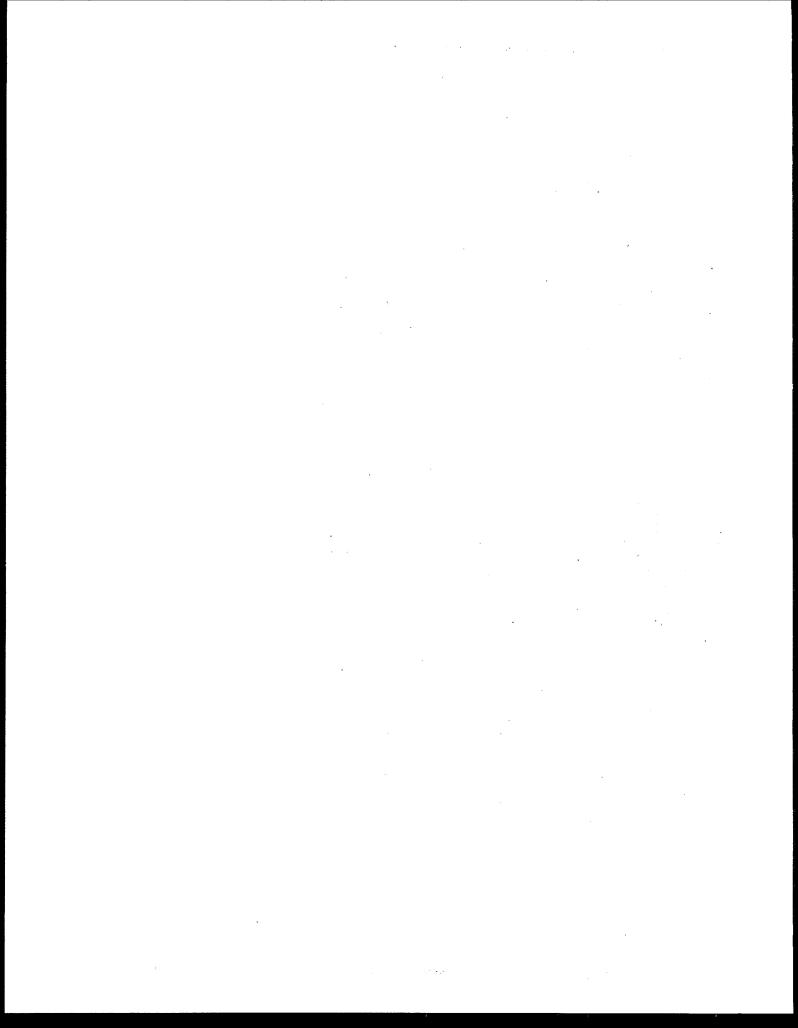


# APPENDIX 6 WASTE CHARACTERIZATIONS FOR TOP 150 ROUTINELY GENERATED COMBUSTED WASTE STREAM COMBINATIONS

EPA reviewed the composition of the 150 largest-quantity wastestream combinations to identify the top 100 containing metals and/or halogenated organics. The following table provides the following information:

- each of the attributes that define a wastestream combination (i.e., RCRA code, SIC code, BRS source codes, BRS form code), and a check mark beside each that played a critical role in EPA's characterization of waste composition;
- the number of wastestreams comprising the combination;
- the number of facilities generating the wastestream combination;
- the constituents and concentrations present;
- the source of information for the characterization of constituents;
- a key noting the type of constituents present
  - "0" denotes that neither metals nor halogenated organics are present
  - "1" denotes the presence of metals
  - "2" denotes the presence of halogenated organics
  - "3" denotes the presence of both metals and halogenated organics;
- principal assumptions used in interpreting the data sources and assigning concentrations.

The information in this appendix is subject to a number of important caveats and limitations, as described in Chapter 2 of the accompanying report.



Appendix 6 Top 150 Routinely Generated Combusted Mastes: Constituent and Concentration Listing

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N	+	•				3.	
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Hathano) Hathyl isobstyl ketone Cyclohasenone Ethyl ester Ethyl ester Ethyl ester Ethyl ester Ethyl ester Acrylonitrile Benzene Phenol		3010	Mathylene chloride Ethylene dichloride Hethyl isobutyl ketone Toluene Chlorobenzene Hydrochloric acid	formaldehyde Malaic anhyd:ide Phthalic anhyd:ide	Acrylic acid Ethyl acryla Acrolain n.Butanul Chromium	Constituents	
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•		•	^			1.07	
ma (c			toluene ant f005 toluene Averat hiorobenzene f1005 documents for f1 documents for f1 dichloride, mett toluene were be concentrations o concentration o epproximately g1 equeous liquids to constituents to a pH of 2 to characteristic.	Assun	they are assumed to concentrations were 1002 characteristic concentrations of or	Assumptions	
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To a			white out to me to	0.00	oun oun	:	
match: ACRA codes and form code.			rio) code assumed due to methy! isob toluene and 6005 code assumed due to thlorobarzene from BDAI background documents for FOO1 FOO5. Concentrat for methylene chloride, ethylene dichloride, methyl isobutyl hatone, tuluene were based on judgement, toluene were based on judgement, toncentrations were adjusted so that concentration of constituents was approximately 10 percent by weight i equeous liquids. Hydrochloric exide to constituents to account for DOO2 characteristic. Concentration of Hi on a pH of 2 to satisfy DOO2	(on-entration of acrying action of contentration of the formation used for the Generator St match. R(RA codes and SIC code.			
			FIGO) code assumed due to methy! Isobuty! Isobut	or si	they are assumed to account for DOO! they are assumed to account for DOO! ton-ent-ations were based on judgment tings characteristic assumed because of concentrations of organic acids in itgui		
	\$ we ve v		FIGO) code assumed due to methyl isobutyl betone and 6005 code assumed due to toluene. Average concentration for thiorobenzene from BDAI background documents for FOGI FOGS. Concentrations for methylene chloride, ethylene dichloride, methyl isobutyl betone, and toluene were based on judgement. Concentrations were adjusted so that total concentration of constituents was concentration of constituents was approximately 10 percent by weight for agenous liquids. Hydrochloric acid added to constituents to account for DOO2 the acteristic. Concentration of Hill based on a pH of 2 to satisfy DOO2	(one entration of errying extended of the setting 1002 characteristic Assumption used for the Generator Survey match. R(RA codes and SIC code.	they are assumed to account for DOO!  (not entrations were based on judgement  (002 characteristic assumed because of low contentrations of organic acids in liquid  one entrations of organic acids in liquid		
			-	1	<u> </u>	_	

Appendix 6
Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RAW		RCRA weste code		SIC Code		Source Code		form Lode	Quantity	3	1.00	Constituents Co	PRC. (P	p=)	Constituent Source	Kay	Assumptions
5							1	B206	64,191	450	450	Xylene Toluene 1,1,1-Trichlorosthane Benzene Tetrachlorosthylene Trichlorotrifluorosthane Naphthalene Trichlorosthylene Oichlorodifluoromethane Benzo(a)anthracene Benzo(a)pyrene Arsenic Barium Cadmium Chromium Lead Zinc	3,	100 200 100 900 800 500 250 20 8 10 65 5	Used O11 RIA	3	All constituents and concentrations were ubtained from lable V 38, "Mean Concentrations of Potentially Hazardous Constituents in Used Oil Burned as On specification fuel," in the used oil RIA.
7		D001 D002 D003 D018 D023 D024 D025 D026		2869		A32	•	8110	64,265	2	1	Cresols Benzene Hydrochloric acid Hydrogen sulfide	1	100 500	RCRA codes RCRA-codes None None	0	0001 characteristic assumed from benzene and hydrogen sulfide was added to cover 0003 characteristic. Concentrations for benzene and cresols are 200 times the regulatory levels. Hydrogen sulfide concentration was based on proposed EPA guidelines. Hydrochloric acid added to constituents to account for D002 characteristic. Concentration of HCl based on a pH of 2 to satisfy 0002 characteristic.
8	,	D001 D018 D035 F003		2869	Ĺ	A37 A57	•	B207 B219	62,350			Acrylic acid Butyl acrylate	50,0 50,0	000	BRS CAS numbers BRS CAS numbers		DOO2 characteristic assumed due to acrylic acid. Concentrations of acrylic acid and butyl acrylate were based on judgement.
9			Ĺ		Ĺ		_		54,912	1	,	Benzene Methyl ethyl ketone	200,0 100,0	000	RCRA waste code RCRA waste code	٥	DOO1 characteristic and FOO3 code assumed due to benzene and methyl athyl ketone. Concentrations were obtained from the Solvents LDR RIA.
,		D001 D002 D003 F002 F020 F024 K017 K018 K020 K028		2821		A37	•	B219	48,039	2		1,1-Dichloroethene Chloroethene 1,2-Dichloroethene 1,1,2-Trichloroethene 1,1,1,2-Tetrachloroethene 1,1,1,2-Tetrachloroethene 1,1,1,2-Tetrachloroethene 1,1,2,2-Tetrachloroethene 1,1,1-Trichloroethene Vinyl chloride Toluene Mathanol Tetrachlorobenzene 2,4,5-Trichlorophenol 2,3,7,8- Tetrachlorod benzo(p) dioxin Hydrogen sulfide	50,0 50,0 50,0 50,0 50,0 50,0 50,0 1,0 1,0	00 00 00 00 00 00 00 00 00 00 00 00			Constituents were obtained from the background document for No18, K020, K028, and F020. F002 code assumed due to 1,1,1 trichloroethane. Vinyl chloride was added to account for K020. Hydrochloric acid added to constituents to account for D002 characteristic and hydrogen suifide added to account for D003 characteristic. Concentration of HCl based on a pH of 2 to satisfy D002 characteristic. Concentration of HCl based on a pH of 2 to satisfy D002 characteristic. Concentration of hydrogen sulfide was based on proposed EPA guidelines. Concentrations of the organics were based on judgement.

Appendix 6

Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RANK	<del></del>	RCRA waste code		SIC ode		ource Cede		or <b>m</b>	Quantity	VS.	fac	Conseitants	Conc. (ppm)	Constituent Source	Kay	Assumptions
10	1	D018 D038	1	2869	-	A34	•	B102	44,977	1	1	Pyridine Ammonia Benzene Toluene (yenide	500 250 250	BRS CAS numbers BRS (AS numbers BRS (AS numbers TC HIA Back IC RIA Back	0	Constituent data taken from Table 3 of "Estimates of Waste Generation by the Griganic Chemical Industry" of the IC LDR RIA Background Document assuming wastewater was derived from the production of pyridine (DO3B). Pyridine concentration was assumed to be 200 times the regulatory lavel. Concentration for ammonia was based on judgement.
11	,	D001 D002		2869		A33		B201	40,001	1	1	Methenol Hydrochloric acid Methyl acetate	100,000 500 10,000	BRS LAS numbers BRS CAS numbers None	0	0001 characteristic assumed from methanol and 0002 characteristic assumed from hydrochloric acid. Heathyl acetate was added since the BRS waste description suggested the presence of organic esters. Concentrations for methanol and methyl acetate were based on judgement. Concentration of HCl based on a pH of 2 to satisfy 0002 characteristic.
12	4	F002	-	2879		A35	7	6101	37,447	;		Methylene chloride	12,000	BRS CAS number	2	Assumed FOO2 code was due to methylene chloride alone. Concentration for methanol was obtained from the Generator Survey. Assumption used for the Generator Survey match: RCRA codes and form code.
13	,	Da01 D018 D035 F003		<b>7869</b>		AIS	,	8207	36,916	1		Benzene Methyl ethyl ketone Xylene Acetone Ethyl ecetete Ethyleneane Ethyl ether Hethyl isobutyl ketone n-Butenol Cyclohexenone Methenol	100,000 30,000 50,000 25,000 25,000 25,000 40,000	RCRA waste code RCRA waste code Solvents LDR RIA Solvents LDR RIA		DOOI characteristic assumed from benzene and methyl ethyl ketone. Average concentrations of benzene and methyl othyl ketone were obtained from the Solvents LDR RIA. All constituents (and concentrations) listed under FOO3 in the Solvents LDR RIA were included in this table. Since the form code indicates a concentrated equeous solution of other organics, concentrations were based on Judgement and their relative proportions as given in the Solvents LDR RIA.
34		D001 D002 D007		2869		A33	,	B602	36,709			1 Chromium Toluene Xylane Lead Cadmium Hydrachloric acid	50,000 50,000 6	RCRA waste code Gen. Survey Gen. Survey Gen. Survey Hone		Concentration for chromium was assumed to be 200 times the regulatory level. Hydrochloric acid added to constituents to account for D002 characteristic. Concentration of HCl based on a pH of 2 to satisfy D002 cheracteristic. Only constituents occurring in more than 33 percent of waste streams with D001 as the only RCMA code (as reported in the Generator Survey) were included to account for the D001 cheracteristic. For these constituents the median concentrations of the wastes in the Generator Survey were used.

Appendix 6
Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RAHK	T	RCRA weste code	T	SIC	Τ.	Source	T	form	Quantity	Γ.	Τ.	I.	T		
<u> </u>	<u> </u>		<u></u>	ode.	<u>L</u>	Code		Code	dasucity	vs	100	Constituents Conc. (ppm	Constituent Source	Key	Assumptions
15		1001 1002 f003 f005 k038 P094		2879		A37	*	B101	35,136	1	1	The triple of triple	D BRS CAS numbers D BRS (AS numbers D F001-F005 Back D Doc. F001-F005 Back D Doc. K03B BDAT Doc.	5	P001 and D002 codes assumed from K038 constituents. P094 is phorate which is one of the constituents of K038. Average concentrations of constituents in spent solvents from production of cyclic esters and phosphoroemidothioate are used for f003 and f005 (oncentration for methylene chloride was based on judgement Hydrochloric acid added to constituents to account for D002 characteristic.  Concentration of HC1 based on a pH of 2 to satisfy D002 characteristic.  Concentrations were adjusted so that total concentration of constituents was approximately 10 percent by weight for aqueous 14quids.
16	4	D001 D005 D006 D007	(	2869	,	A 33	7	8219	32,779	18	17	[marnanor 50,000	BRS CAS numbers BRS CAS numbers BRS CAS numbers	0	D001 characteristic assumed due to the organics. Concentrations for organics from the Generator Survey. Assumption used for the Generator Survey match: exact match on all data elements.
		D008 D018 D026 D035 F001 F002					,	6719	31,340	1	1	Sarium   20,000	RCRA waste code Solvents LDR RIA	3	D001 characteristic assumed to be due to the organics denoted by the D codes. Concentrations of constituents indicated by the D codes were assumed to be 200 times the regulatory levels. All constituents included in the LDR RIA for F001 and F002 were added. Since the form code indicates unspecified organic liquids, concentrations were based on judgement and their relative proportions as given in the Solvents LDR RIA.
16		F002 F005		2834		A37		B101	28,640	1		Methylene chloride         1,000           Toluene         1,000           Xylene         1,000           Acetone         10           Methanol         1	BRS CAS numbers BRS CAS numbers BRS CAS numbers BRS CAS numbers BRS CAS numbers		F002 and F005 codes were assumed to be due to organics indicated by the BRS CAS numbers. Concentration for xylene was based on judgement. Concentrations for other constituents obtained from the Generator Survey. Assumption used for the Generator Survey match: RCRA codes and form code.
				,		nJ/		B102	27,247	1		5,000     Dichloropropens   200     Dichloropropens   200     Chloroform   200     Mathyl chloride   30     Carbon tetrachloride   30     Tetrachloroethane   30     Mathylene chloride   30     Selenium   30			Constituent information used for waste stream in the Generator Survey with the RCRA codes D001-D003 and the SIC code 2818 (Organic Pasticide Products). Hydrochloric acid added to constituents to account for D002 characteristic. Concentration of HCl based on a pH of 2 to satisfy D002 characteristic. Concentrations were adjusted so that total concentration of constituents was approximately 10 percent by weight for equeous liquids.

Appendix 6
Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RAMK		RCRA was	le code		SIC ode	ource Code	orns ode	Quantity	8	fac	Constituents Conc.	(pp=)	Constituent Source	Key	Assumptions
20	1	D001 D018	(1019 (1039		2669	A74	8202	26,708	1	1	1.1.2 Irichloroethane 1.1.2.2 fetrachloroethane Hexachloro 1.3 butadiene Benzane Carbon tetrachloride Tetrachloroethylene 1.2-Oichloroethane 1.1-Dichloroethane 3-Chloropropene Trans 1.3-dichloropropene 1.2-Oichloropropene 2-Chloro-1.3-butadiene Cis 1.3-dichloropropene	500 500 60 1,500 1,500 30,000 15,000 15,000 9,000 9,000 9,000 4,000	BRS CAS numbers BRS CAS numbers BRS CAS numbers RCRA code RCRA code RCRA code BDAT Back Doc	2	Based on the BRS description, the waste was assumed to be 90 percent mixed heavy ends generated during the production of chlorinated hydrocarbons. The remaining 10 percent comprised of toxics are listed hare. Concentrations represent midpoints of ranges indicated for constituents in untreated f024. The concentrations were adjusted so that the total organic concentration was approximately 10 percent
- 21	,	D001 F002	F003 F005	1	2833	A35	B101	26,284	į	1	Methanol Acatone Toluene Hathylene chloride	30,000 20,000	BRS CAS numbers BRS CAS numbers BRS CAS numbers BRS CAS numbers	2	F002, F003, and F005 codes were assumed to be due to organics indicated by the BRS CAS numbers. Constituent concentration obtained from the Generator Survey and were adjusted so that total concentration of constituents was approximately 10 percent by weight for aqueous liquids. Assumption used for the Generator Survey match. RCRA codes and SIC code.
22		0001 D002	D003 D018		.'Bb's	A 35	8219	25,847	. 1	1		40,000 500	RCRA codes RCRA codes None None	0	D001 characteristic assumed from benzene and hydrogen sulfide was added to cover 0003 characteristic. Concentration for benzene based on judgement, concentration for cresols was assumed to be 200 times the regulatory level, and concentration for hydrogen sulfide was based on proposed EPA guidalines. Hydrochloric acid added to constituents to account for D002 characteristic. Concentration of HCl based on a pH of 2 to satisfy D002 characteristic.
23	,	0001						25,371	223	178	Ethanol Epichlorohydrin 1,2-Dichloropropane 1,2,3-Trichloropropane Bis (2-chloroethyl) ather	50,000 10,000 10,000	BRS waste desc. BRS waste desc. KO17 listing KO17 listing KO17 listing	2	Based on the BRS waste description, the waste was assumed to be a KO17 weste (heavy ends from the purification column in the production of epichlorohydrin); all constituents for which KO17 nonwastewaters is listed were added. Constituent concentrations were based on judgement.

Appendix 6
Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RANK	RLRA weste code		SIC Ioda	ource Cede		orm ode	Quantity	8	inc	Constituents	Conc. (ppm)	Constituent Source	Key	Assumptions
1	K022	,	2849	A33	•	8606	23,281		1	Phenol Acenaphthalene Acenaphthene Anthracene Diphenyl mine Fluorenthene Naphthalene Ethylbenzene Acetophenone Acetone Aniline Styrene Toluene Rylene Fluorene Antimony Arsenic Berfum Beryllium Cadmium Chromium Copper Lead Mercury Nickel Silver Thallium Vanadium Zinc	10,000 10,000 10,000 10,000 10,000 10,000 10,000 2,500 100 15 100 15 100 15 0.01 0.01 0.01	BRS (As number Gen. Survey	1	ioncentrations of all constituents were obtained from the Generator Survey. Assumption used for the Generator Survey match. RiHA codes and SIC code. Concentrations of the organics were adjusted to account for the form code that indicates organic sludges.
25	1001 (2004 (2005 (2006 (2007					B204	22,251	3	3	Barium Arsenic Chromium Cadmium Toluene Xylene Lead	1,000 1,000 200 50,000 50,000	RCRA codes RCRA codes RCRA codes RCRA codes Gen. Survey Gen. Survey		It was assumed that this waste stream combination is a mixture of DOOI wastes and other wastes. Only constituents occurring in more then 33 percent of waste streams with DOOI as the only RCRA code (as reported in the Generator Survey) were included to account for the DOOI characteristic. For these constituents the median concentrations of the wastes in the Generator Survey were used. Concentrations for metals were assumed to be 200 times the regulatory levels. Assumed that no halogenated organic constituents were present in this waste although the form code indicates the potential presence of halogenated organics.

Appendix 6

Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RAHK	Ī	RCRA waste code	IC ode		ource Code		or <b>m</b>	Quantity	45	f ac		-)	Constituent Source	Kay	Assumptions
26		<b>▶048</b>	2911	1	A75		8503	19,996	1	1	Naphthalene Anthracene Phenanthrene Benzo(a)pyrene Benzo(a)pyrene Bis (2-ethylhexyl) phthalate Dimethyl phthalate Ethylbenzene Phenol Chrysene Cresols Toluene Pyrene Antimony Arsenic Berlum Beryllium Cadmium Chromium Copper Lead Mercury Nickel Selenium Silver Thallium C	50 20 50 20 5 5 20 5 5 20 5 5 20 5 5 20 5 5 20 5 5 20 5 5 20 5 5 20 5 5 20 5 1 20 5 1 20 1 20	Gen. Survey	3	Assumption used for the Generator Survey match. RERA codes, SEC code, and source code.
21	,	(1001 D002 D003 D008 11018 D023 D024 D025 D026	2869		Et A	•	B219	18,825	2	1	Benzene 50,0 Lead 1,0 Hydrogen sulfide	000 000 500	RCRA waste code RCRA waste code RCRA waste code None None	1	D001 characteristic assumed due to benzene Hydrogen sulfide included to account for D003 characteristic. Concentration of hydrogen sulfide based on proposed EPA guidelines. Hydrochloric acid added to constituents to account for D002 characteristic. Concentration of HCl based on a pH of 2 to satisfy D002 characteristic. Toncentration for lead was assumed to be 200 times the regulatory level and concentrations for benzene and cresols were based on judgement.
28	ľ	D001	2869		A33	,	B203	18,774	11	3.1	Toluene 50,6 Butanol 125,6 Propanol 50,6	000 000 000	BRS CAS numbers BRS CAS numbers BRS waste desc. BRS waste desc. BRS waste desc.	0	DOD1 characteristic assumed due to organics. Concentrations for propanol and morpholine were based on judgement. Concentrations for the other constituents were obtained from the Generator Survey. Assumption used for the Generator Survey match: exect match on all data elements.

Appendix 6
Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RAM		PK.	RA weste	code		SIC Code	ource Code		ole ode	Quantity	3.	100	Constituents	Conc. (ppm)	Constituent Source	Key	Assumptions
29	,		01 Ł005 ŁN			2384	A37	•	8101	18,747	1	1	Hethanol Acetone Hathylane chloride Hethyl isobutyl ketone n-Butanol Toluene Hathyl ethyl ketone Banzene	25,000 8,000 4,000 1,500 30,000 2,000	BRS CAS numbers BRS CAS numbers BRS CAS numbers BRS CAS numbers BRS CAS numbers GRS CAS numbers Gen. Survey Gen. Survey	5	DOD1, FUD2, and FOD3 codes are assumed from organics indicated by the BRS CAS numbers. Banzene, methyl athyl ketone, and toluene were obtained from the Generator Survey and used to account for FOD5 code. Concentrations of all constituents were obtained from the Generator Survey. Assumption used for the Generator Survey match: RCHA codes and source code. Concentrations were adjusted so that total concentration of constituents was approximately 10 percent by weight for aqueous liquids.
30	,		01 0003 00			2869	A35	•	8212	18,603	1	1	Benzene Hydrocyenic ecid	100,000 100	RCRA waste code HCRA waste code	0	DOO! characteristic assumed due to benzene and DOO3 characteristic assumed due to hydrocyanic acid. Concentration for hydrocyanic acid obtained from PO63 background document. Benzene concentration based on judgement.
		D00	)1 F002 F0		1	2831	A 35	•	<b>0</b> 103	18,154	J	1	Hethylene chloride Hethenol Acetone	10,000	BRS CAS numbers BRS CAS numbers BRS CAS numbers		All the D and F codes are assumed from organics indicated by the BRS CAS numbers. Concentrations were obtained from the Generator Survey and adjusted so that total concentrations of constituents was approximately 10 percent by weight for aqueous liquids. Assumption used for the Generator Survey match: RCRA codes and SIC code.
1.	•	1410				2869	A 3 3	*	8206	17,367	2	2	Toluene Methenol	30,000 50,000	BRS CAS numbers BRS CAS numbers	0	DOO1 characteristic assumed due to organics indicated by the CAS numbers. Concentrations were obtained from the Generator Survey. Assumption used for the Generator Survey match: RCRA codes and form code.
	,	,	8 DO 38 KO	22 K083		2865	<b>A</b> .		В .	17,303	1	1	Pyridine Benzene Acetophenone Phenoi Aniline Diphenylamine Mitrobenzene Mickel	10 250,000 50,000 10 10	RCRA waste code RCRA waste code KO22 back. doc. KO22 back. doc. KO83 back. doc. KO83 back. doc. KO83 back. doc. KO83 back. doc.		Concentration of benzene and KO22 constituents obtained from KO22 8DAT background document; concentrations for KO83 constituents obtained from KO83 BDAT background document; Pyridine concentration assumed to be 200 times the regulatory lavel. Nonwastewater form was assumed based on descriptions of KO22 and KO83.

Appendix 6

Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RANK	Γ	RCRA w	aste code	I(.	ource Code		orm ode	Quantity	Vs.	101	Constituents	Conc. (ppm)	Constituent Source	Key	Assumptions
34	1	f003 f0	05		·		B219	17,218	12		Methy) ethyl ketone Acetone Toluene Ethyl acetete Xylene n-Butenol Methyl isobutyl ketone Ethylbenzene Chlorobenzene Arsenic Barium Chromium Lead Silver Cadmium Selenium Mercury	125,000 100,000 90,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000	Gen. Survey		Assumption used for the Generator Survey match HiHA codes and form code (oncentration for barium was adjusted since the Generator Survey indicated an unrealistically high concentration for barium
35		0001 D0 F005	38 FQQ2 FQQ3	2834	A 37		B203	16,380	1	3	Methano'l Acetone Acetonitrile Toluene Xylene Pyridine	350,000 120,000 120,000 750	BRS CAS numbers BRS CAS numbers BRS CAS numbers BRS CAS numbers BRS CAS numbers BRS CAS numbers RCRA waste code	0	All the D and F codes are assumed from organics indicated by the BRS CAS numbers. Concentrations were obtained from the background document for F001 F005 for the pharmaceuticals industry. The waste code F002 and form code B203 conflict since F002 indicates halogenated organics while B203 indicates only non-halogenated organics Considering the constituents indicated by the BRS CAS numbers, the form code was assumed to be more accurate than the waste code. Since the form code indicates a solvent mixture, concentrations were adjusted so that total concentration would be approximately 1,000,000 ppm.
36	,	D001 D0	DO2 DO07	2869	A33	,	8219	16,099	2	2	Chromium Vinyl scetate Toluene Xylene Lead Cadmium Hydrochloric acid	50,000 50,000 50,000 50,00	RCRA weste code BRS weste desc. Gen. Survey Gen. Survey Gen. Survey None	1	Concentration for vinyl acetate was based on judgement. Only constituents occurring in more than 33 percent of waste streams with D001 as the only RCRA code (as reported in the Generator Survey) were included to account for the D001 characteristic. For these constituents the median concentrations of the wastes in the Generator Survey were used. Concentration of chromium was assumed to be 200 times the regulatory level. Hydrochloric acid added to constituents to account for D002 characteristic. Concentration of HCl based on a pH of 2 to satisfy D002 characteristic.

Appendix 6
Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RANK		RKRA meste code	SIC Code	Source Code	form Code	Quantity	100	) ac	Constituents	Conc. (ppm)	Constituent Source	Key	Assumptions
37	4	D001 B002	2819	A37	/ B219	15,997	1		Hathanol Hathyl isobutyl ketone Acrylonitrile Acetone Benzene Cyclohexanone Ethyl actate Ethyl ether Ethylene dichloride Benzel chloride Benzel chloride Antline Benzel chloride Formaldehyde Hexachlorocyclopentadie 1,4-Diethylene oxide Phenol Tetrachlorosthylene Toluene Antimony Arsenic Barlum Beryllium Cadmium Copper Lead Mercury Nickel Selenium Silver Thallium Hydrochloric acid	30,000 30,000 30,000 30,000 30,000 30,000 30,000 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500 1,500	Gen Survey Gen. Survey	3	1001 characteristic assumed due to the organics. Hydrochloric acid added to constituents to account for DOD2 characteristic. Concentration of HC1 based on a pil of 2 to satisfy DOD2 characteristic. Assumption used for the Generator Survey match: RCRA codes and form code.
38		D001 D018 D043 F001 F002 F003 F004 F005			√ B204	15,509	1		Benzene Vinyl chloride Cresols Nitrobenzene Tetrachloroethylene Trichloroethylene Methylene chloride 1,1,1-Trichloroethane Carbon tetrachloride Chlorobenzene 1,2-Dichloroethane Trichloroefluoromethane Xylene Acetone Ethyl acetate Ethylbenzene Ethyl isobutyl ketone n-Butenol Cyclohexanone Methanol	100 400 100,000 50,000 50,000 50,000 40,000 50,000 30,000 50,000 50,000 50,000 50,000 25,000 25,000 25,000	RCRA waste code RCRA waste code Solvents LDR RIA		D001 characteristic and f005 code assumed due to benzene. All constituents included in the IDR RIA were added. Concentrations for benzene and vinyl chloride were assumed to be 200 times the regulatory level. Since the form code indicates a solvent mixture, concentrations were adjusted so that total concentration would be approximately 1,000,000 ppm.

Appendix 6

Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RANK	Γ	RCRA waste code	JC ode	our ce Ceda		orm ode	Quantity	VS	# Fac	Constituents	Conc. (ppm)	Constituent Source	Key	Assumptions
39		D001 F024	2819	A33	,	B219	14,893	1	1	1,3-Dichloropropylene 1,2-Dichloropropane Allyl chloride 1,2,3 Trichloropropane Acetone Benzene	150,000 100,000 50,000 10,000	BAS CAS numbers BAS CAS numbers BAS CAS numbers Gen. Survey F024 BDAT Doc. F024 BDAT Doc.	2	tuncentrations for 1,3-dichloropropylene and 1,2,3 trichloropropane were obtained from the Generator Survey. Acetone and benzene were added from the F024 background document to account for the D001 characteristic Concentrations for the other constituents are from the F024 Background Document, and ere the midpoints of the possible concentration range for each constituent. Assumption used for the lenerator Survey match: RCRA codes and form code.
40	,	D018 K022	2821	A33	,	B208	14,636	1	1	Benzene Acetophenone Phenol	750,000	BRS CAS numbers BDAT Back. Doc. BDAT Back. Doc.	0	Acetophenone and phenol concentrations were obtained from the K022 Background Document, and are the midpoints of the possible concentration range for these constituents, benzene concentration was based on judgement. The concentrations of the phenolics were adjusted since the form code indicates a concentrated phenolic liquid
41	•	D001 F001 F002 F003 F005 U001 U002 U003 U019 U028	2869	A33		8219	14,217		1	Acetaldehyde Acetone Acetonitrile Benzene Disthylhexyl phthalate Tetrachloroethylene Irichloroethylene Mathylene chloride 1,1,1-Trichloroethene Carbon tetrachloride Chlorobenzene 1,2-Dichlorobenzene Irichlorofluoromethene	50,000 50,000 50,000 50,000 50,000 50,000 50,000 50,000 50,000	RCRA waste code RCRA waste code RCRA waste code RCRA waste code RCRA waste code Solvents LDR RIA Solvents LDR RIA	2	DOOI characteristic assumed due to organics. F003 code assumed due to acetone and F005 code assumed due to benzene. All constituents included in the Solvents LDR RIA for F001 and F002 were added Concentrations for all constituents were based on judgement.
42		U001 D002 D003 D018 D026 D035 F002 F003 F004 F005	2869	A33 .		B219	14,194	2	1	Cresols Methyl ethyl ketone Benzene Tetrachloroethylene Methylene chloride Trichloroethylene Chlorobenzene 1,1,1-Trichloroethene 1,2-Dichlorobenzene Trichlorofluoromethene Hydrogen sulfide Hydrochloric acid	50,000 50,000 50,000 50,000 50,000 50,000 50,000 50,000	RCRA waste code RCRA waste code RCRA waste code Solvents LDR RIA Solvents LDR RIA Solvents LDR RIA Solvents LDR RIA Solvents LDR RIA Solvents LDR RIA Solvents LDR RIA None Hone	2	DOO! characteristic assumed due to organics, FOO3 assumed due to methyl ethyl ketone, FOO4 assumed due to cresols, and FOO5 assumed due to benzene. Hydrochloric acid added to constituents to account for DOO2 characteristic. Concentration of HCl besed on a pH of 2 to satisfy DOO2 characteristic. Hydrogen sulfide was added to account for DOO3 characteristic. The concentration for hydrogen sulfide was based on proposed EPA guidelines. All constituents included in the Solvents LDR RIA for FOO2 were added. Concentrations for organic constituents were based on judgment.
43	1	D002	2833	A34		8105	13,462	ı	1	Sulfuric acid	500	BRS CAS numbers	0	Concentration of sulfuric acid based on a pH of 2 to satisfy DOO2 characteristic.

Appendix 6
Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RAHK		RCRA weste code		SIC ode	ource Code		orm ode	Quantity	<b>%</b>	Fac	Constituents Co	nc. (pjm)	Constituent Source	Key	Assumptions
44		D001 D00R		2621	A33		B602	13,195	3		Methyl methacrylate lead	90,000 4,000	BRS (AS numbers RCRA waste code	1	INDD1 characteristic assumed from methyl methacrylate. Concentration for methyl methacrylate was obtained from the Generator Survey. Concentration for lead was adjusted since the Generator Survey indicated an unrealistically high concentration for lead. Assumption used for the Generator Survey match: RCRA codes and form code
45	,	D001 D002 D003 D004 D005 D006 D007 D008 D009 D010		2869	A33		8105	13,182	1	1	Arsenic Barium Cadmium Chromium lead Marcury Selenium Toluene Hydrochloric acid Hydrogen sulfide	5 0.5 0.05 0.5 50,000 50,000	RCHA waste code RCRA waste code RCRA waste code RCRA waste code RCRA waste code RCRA waste code RCRA waste code Gen., Survey Gen. Survey None	1	Only constituents occurring in more than 33 percent of waste streams with D001 as the only RCRA code (as reported in the Generator Survey) were included to account for the D001 characteristic. For the organics the median concentrations of the wastes in the Generator Survey were used. Hydrochloric acid added to constituents to account for D002 characteristic and hydrogen sulfide added to account for D003 characteristic. Concentration of HC1 based on a pH of 2 to satisfy D002 characteristic. Concentrations of hydrogen sulfide were based on proposed EPA guidelines. Concentrations for metals were obtained from the Generator Survey. Assumption used for the Generator Survey match: RCRA codes.
46	,	F017 K019 K020		2869	A33	•	8601	13,073	1	1	Ethylene dichloride 1,1,2-Trichloroethane Tetrachloroethane Tetrachloroethane Trichloroethane Trichloroethylene 1-chloro, 2,3-epoxy propan	50,000 50,000 50,000 5,000	Gen. Survey Gen. Survey Gen. Survey Gen. Survey Gen. Survey Gen. Survey	2	Assumption used for the Generator Survey match: RCRA codes and form code. Concentrations for ethylene dichloride and 1,1,2-trichloroethane were adjusted since the Generator Survey indicated unrealistically high concentrations.
47		D001	•	2869	A35	,	8219	12,842	6	6	Methanol Isobutanol Acetone n-Butanol Methyl ethyl ketone Arsenic Barium Lead Chromium Cadmium Mercury Selenium Silver	50,000 750 50 0.05 0.05 0.5 0.5 0.5 0.05	Gen. Survey Gen. Survey Gen. Survey Gen. Survey	1	Assumption used for the Generator Survey match: exact match on all data elements.
48	′	D001 F003	,	2833	A33	′	<b>8</b> 203	11,772	2	2	Acetone Hethanol		BRS CAS numbers BRS CAS numbers	0	DOD1 characteristic and F003 code were assumed to be due to organics indicated by the BRS CAS numbers. Constituent concentrations obtained from the Generator Survey. Assumption used for the Generator Survey metch: RCRA codes and form cude.

Appendix 6

Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RAHK		RCRA weste code		SIC ode	ource Code		ormi ode	Quantity	45	100	Constituents Conc. (ppm)	Constituent Source	Key	Assumptions
49	1	k027		2865	A33	*	B4U9	11,123	1	1	Toluene dissorvanate 10,000 2,4 Toluene diamine 100,000 2,6 Toluene diamine 100,000	Gen Survey Gen Survey Gen. Survey Gen. Survey Gen. Survey	3	Assumption used for the Generator Survey match: RCRA codes. Concentrations for the organics were adjusted to account for the form code that indicates organic solids
50	,	0001		2869	A37	,	8207	11,114	4	3	Ethyl acrylate 35,000 Acrylic acid 35,000	BRS CAS numbers BRS CAS numbers BRS CAS numbers BRS CAS numbers	O	Acrylic acid assumed to account for DOOI characteristic. Concentration of sulfuric acid based on a pH of 3 since this waste stream combination does not exhibit the DOO2 characteristic. Concentrations of other constituents based on judgement.
51	,	D001 D002 F003	,	2834	A31	1	8101	10,976	1	1	Cyclohexane 25,000 Isopropanol 25,000 Acatona 25,000 Acatona 600	BRS CAS numbers BRS CAS numbers BRS waste desc. BRS waste desc. BRS waste desc. BRS waste desc.	0	DOO) characteristic assumed due to the organics, DOO2 characteristic assumed due to acatic acid and HCl, and FOO3 code assumed due to methanol. Constituent concentrations were based on judgement. Concentrations of acatic acid and HCl based on a pH of 2 to satisfy DOO2 characteristic.
52	,	D001 F001 F002 F005				,	B204	10,929	2	2	Acrylonitrile 125,000 Mathyl chloroform . 75,000 Acatone 10,000 Chlorobenzene 1,000 Cyclohexanone 1,000 4,4-mathylene bis(2-chloroanillre0) Mathyl enbyl ketone 10,000 Mathyl sthyl ketone 10,000 Tetrachloroathylene 50,000 1,1,2-Trichloroathane 75,000 Trichloroathylene 35,000	Gen. Survey	2	DOOl characteristic assumed due to organics. Assumption used for the Generator Survey match: RCRA codes.
53	,	D001 D005 D006 D007 D008		7 389	A89		B204	10,883	1		Cadmium   200   Chromium   1,000   Lead   1,000   Lead   1,000   Mathyl ethyl ketone   50,000   Mathyl isobutyl ketone   50,000   n-Butanol   50,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000   1,000	D RCRA waste code D RCRA waste code RCRA waste code D RCRA waste code Gen. Survey D Gen. Survey D Gen. Survey D Gen. Survey	1	Constituents from the Generator Survey that could account for the DOOI characteristic were added. Concentrations for metals were assumed to be 200 times the regulatory level. Concentrations for organics were based on Judgement. Assumption used for the Generator Survey match: RERA codes. Assumed that no halogenated organic constituents were present in this weste although the form code indicates the potential presence of halogenated organics.

Appendix 6
Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

KANK		RERA waste code	SIC Code	Source Code	form Lode	Quant ity	3	100	Constituents Con	nc. (ppm)	Constituent Source	Key	Assumptions
54		F022	/ 2869	A33	1 B<19	10,846	1		thylbenzens Acetophenons Phenol Acetone Acenaphthelens Acenaphthene Aniline Anthracene Diphenyl amine Fluorenthene Fluorene Naphthalene Styrene Toluene Hylene Antimony Arsenic Barium Beryllium Cadmium Chromium Cropper Lead Mercury Mickel Selenium Silver Thallium Vanadium Zinc	40,000 30,000 25,000 100 100 100 11 500 15 0.01 0.01 0.01	Gen. Survey		Assumption used for the Generator Survey match. HIRA codes and SIC code.
56		14001 F901 F002 F003	2869	A35	/ 8204	10,782	8		Acetone Hethylene chloride 1,1,1-Trichloroethene Hethyl chloroform Trichlorotrifluoroethene Selenium	500,000 220,000 60,000 60,000 60,000	Gen. Survey Gen. Survey Gen. Survey Gen. Survey Gen. Survey Gen. Survey	İ	0001 characteristic assumed due to organics. Assumption used for the Generator Survey match: RCRA codes and form code. Since the form code indicates a solvent mixture, concentrations were adjusted so that total concentration would be approximately 1,000,000 ppm.
57	-	D018 F037 F038 K048	2911	A75	/ B603	10,732	2		Toluene Kylene Lead Chromium Cadmium	50,000 5 5 0.1	Gen. Survey Gen. Survey Gen. Survey Gen. Survey Gen. Survey	1	Only constituents occurring in more than 3: percent of waste streams with D001 as the only RCRA code (as reported in the Generator Survey) were included to account for the D001 characteristic. For these constituents the median concentrations of the wastes in the Generator Survey were used. Concentrations of the organics were adjusted to account for the form code that indicates organic liquids.
		KO49 KO50 KO51			6003	10,580	1	- 1	Benzene Lead Chromium	1,000	KO48-51 List.	1	Benzene assumed to account for all RCRA codes; lead and chromium were added based upon the 8040-5; listing document. Concentrations of the metals were assumed to be 200 times the regulatory level. Benzene concentration was based on judgement.

Appendix 6

Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

9.00%	Γ	RCRA waste code	il(	ource Code		orm ode	Quantity	Š	) toc	Constituents	Conc. (ppm)	Constituent Source	Key	Assumptions
58	7	k022	2065	A13		8602	9,432	1	1	Acatona Phanol Arsanic Cadmium Chromium	50,000 5 0 5	BRS waste desc. Gen. Survey Gen. Survey Gen. Survey Gen. Survey	1	Concentration for acetone was based on Judgement Concentrations for the other constituents were obtained from the Generator Survey. Assumption used for the Generator Survey match: RCRA codes and form code.
59	,	UU01 D011 D018 D021 D022	3861	A49	-	8204	9,390	2	1	Methanol Acatone Toluene Xylene Mathylene chloride Chlorobenzene Chloroform Benzene Silver	150,000 150,000 150,000 150,000 20,000 20,000 20,000	BAS (AS numbers BRS (AS numbers BRS (AS numbers BRS (AS numbers BRS (AS numbers BRS (AS numbers BRCRA waste code RCRA waste code RCRA waste code RCRA waste code	3	DD01 characteristic assumed due to organics. Concentration for silver was assumed to be 200 times the regulatory level. Concentration for the other constituents were based on judgement with the consideration that the BRS waste description indicated that the weste is a bulk solvent waste from chemical manufacturing.
60		D001 F003 F005				B204	8,747	15	15	Acetone Methyl ethyl ketone Toluene Acetonitrile Benzene n-Butenol (hloroform Cyclohaxanone Dichlorobenzene Ethyl ether Methanol Methyl chloroform Methyl chloroform Methyl isobutyl ketone Pyridine Styrene Tetrachloroethylene Trichloroethylene Trichloroethylene Tylone Antimony Arsemic Berlum Beryllium Cadmium Chromium Copper Lead Mercury Nickel Selenium Silver Thallium Vanadium Zinc	100 000 100 000 100 000 10 000 3 000 3 000 10 000 40 000 65 000 4 000 60 000 65 50 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000 60 000	Gen. Survey	3	DOOI characteristic assumed due to organics. Assumption used for the Generator Survey match: RCRA codes and form code.
61	•	D001 D007 D008 D018	2911	A89		B204	8,564			Benzene Chromium Lead	1,000	RCRA waste code RCRA waste code RCRA waste code	1	D001 characteristic assumed due to benzene Concentrations for chromium and lead were assumed to be 200 times the regulatory level. Benzene concentration was based on judgement. Assumed that no halogeneted organic constituents were present in this waste although the form code indicates the potential presence of halogenated organics.

Appendix 6
Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RAHK		HIRA weste cude	I	SIC Code		Source Cede		lorm Lode	Quantity	vs.	fac	Constituents	Conc (ppm)	Constituent Source	Key	Assumptions
62	,	0059 0001 0018 0019 0055		2641		A37	,	8202	8,416	1	1	Vinyl chluride Trichloruethylene Tetrachloruethylene Benzene Carbon tetrachloride Chloroform 1,2-Dichloroethane	100,000 100,000 100,000 100,000	BRS (AS numbers BRS (AS numbers BRS (AS numbers RCRA waste code RCRA waste code RCRA waste code RCRA waste code	2	DOOI characteristic assumed due to organics. Constituent concentrations were based on judgement, for this waste stream combination, the concentrations were not assumed to be 200 times the regulatory level, since the total concentration with such an assumption was less than 3 percent which appears to be too low for organic liquids (as is indicated by the form code)
63		D001 D002 F003 F005				A37		8101	8,326	1	1	Methanol Acatona Ethylane glycol Toluana Hydrachloric acid	25,000 25,000 10,000	BRS CAS numbers BRS CAS numbers BRS CAS numbers BRS CAS numbers BRS CAS numbers None	-	All D and F codes were assumed to be due to organics indicated by the BRS CAS numbers. Constituent concentrations based on two waste streams in the background document for F001 F005, pharmaceuticals manufacturing, except for ethylene glycol, which was based on judgement. Hydrochloric acid added to constituents to account for D002 characteristic. Concentration of HCl based on a pH of 2 to satisfy D002 characteristic. Constituent concentrations were adjusted so that total constituent concentration was approximately 10 percent by weight for aqueous liquids.
65		(AUU 2		2821		A 37	•	#:U6	8,274	2	2	Methanol Acetaldehyde Vinyl acetate Xylene Methyl methacrylate	5,000 5,000 5,000	BRS CAS numbers BRS CAS numbers BRS CAS numbers BRS CAS numbers BRS CAS numbers		D001 characteristic assumed due to organics. Concentration for mathyl methacrylate obtained from the Generator Survey. Concentrations of all other constituents were based on judgement. Assumption used for the Generator Survey match: exact match on all data elements.
66		D004 F002 F005		2869	,	A09		B207	8,001	1	1	1,4-Diethylene oxide Hydrochloric acid Chromium Copper Nickel	3,100 300 10	Gen. Survey Third Third LDR RIA	ì	Given the form code B207 (concentrated aquaous solution of other organics) in combination with the source code A09 (clean out of process equipment), information in the Third Third LDR RIA for D002 wastes with the waste form "spent acid with metals" was used. The only constituent provided by the Generator Survey was added Assumption used for the Generator Survey match: exact match on all data elements. Concentrations of the organics were adjusted to account for the form code that indicates organic liquids.
		DOON FOOL FOOLS		2879		A09	,	B102	7,954	1	1	Ketocarbamate Toluane Xylene Arsenic	25,000 5,000	BRS CAS numbers BRS CAS numbers BRS CAS numbers RCRA waste code	0	F005 code assumed due to toluene. Organic constituent concentration were taken from the F001-F005 background document for the production of n-alkyl carbamate and ketocarbamate, and were adjusted so that total concentration of constituents was approximately 10 percent by weight for aqueous liquids. Arsenic concentration was assumed to be 200 times the regulatory level. Since the BRS CAS numbers did not indicate the presence of any F002 constituents, it was assumed that F002 code was an error.

Appendix 6

Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RANK	_	RCRA weste code	T		IC ede		ource Code		orm obe	Quantity	vs	fac	Constituents	Conc.	(pp=)	Constituent Source	Kay	Assumptions
67				-	3221	*	A64		B206	7,914	2	2	Xylene Toluene Benzene Tatrachloroethylene Benzo(a)pyrene 1,1,1-Trichloroethene Trichloroethylene Arsenic Barium Cadmium Chromium Lead Zinc		200 10 10 10 6 5 5 100 1 8	Used OI1 RIA	3	All constituents and concentrations were obtained from Table IV-15, "Concentration of Potentially Hazardous Constituents in Automotive Used Oil Samples Taken Directly from Generators," in the used oil RIA.
68	,	D001 D005 D006 E	0007				,	,	8407	7,826	2	2	Barium Cadmium Chromium Lead Kylene n-Butanol Mathyl isobutyl ketone		200 1,000 1,000 1,000	RCRA waste code RCRA waste code RCRA waste code RCRA waste code Gen. Survey Gen. Survey	1	Constituents from the Generator Survey that could account for the DOD1 characteristic were added. Assumption used for the Generator Survey match: RCRA codes. Concentrations for metals were assumed to be 200 times the regulatory level. Concentrations for organics were based on judgement since 1) the Generator Survey indicated very low concentrations and 2) the form code indicates an organic solid.
69	,	D001 D002 D003 (D021 D023 D024 (D026 D035			2869		A33	,	8219	7,418	2	1	Methyl ethyl ketone Cresols Chlorobenzene Benzene Hydrogen sulfide Hydrogen sulfide Hydrochloric ecid		40,000 20,000 100 500	RCRA waste code RCRA waste code RCRA waste code RCRA weste code None	2	DOOI characteristic assumed due to organics. Hydrogen sulfide included to account for DOO3 characteristic. Concentration of hydrogen sulfide based on proposed EPA guidelines. Hydrochloric acid added to constituents to account for DOO2 characteristic. Concentration of HCl based on a pH of 2 to satisfy DOO2 characteristic. Concentrations of other constituents were assumed to be 200 times the regulatory level.

Appendix 6

Fop 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

ЯАНК	<u> </u>	RCRA waste code	SIC SIC	Source Code	form Lode	Quantity	3	,,,	Constituents	Conc. (pp	-)	Constituent Source	Kay	Assumptions
70		pun 1002	2049	A33	/ 8219	7,412	2		Methanol Methyl isobutyl kelone Acetone Acrylonitrile Benzene (yclohexanone Ethyl acetate Ethyl acetate Ethyl endichloride Benzel chloride Benzel chloride Chloroform Phenol Tetrachloroethylene Joluene Formaldehyde Carbon tetrachloride J. 4-Diethylene oxide Hexachlorocyclopentedien Antimony Arsenic Berium Beryllium (admium (hromium (hopper Lead Mercury Nickel Selenium Silver Thallium Vanadium Zinc Hydrochloric acid	0.000000000000000000000000000000000000	000 000 000 000 000 000 000 000 000 00	Gen. Survey	3	Assumption used for the Generator Survey match. HiRA codes and form code. Hydrochloric acid added to constituents to account for DOO2 characteristic. Concents ation of HCl based on a pH of 2 to satisfy NOO2 characteristic.
72		D001 D007 D008 D018 D022 D026 D027 D028 D033 D036 D001 D018 D038 F003	2869	A99	<b>√</b> 8219	7,001	1		Cresols 1,4-Dichlorobenzene Chloroform Nitrobenzene Benzene 1,2-Dichloroethene Hexachloro 1,3-butadiene Chromium Lead	15,00 15,00 15,00 15,00 15,00	00 E	RCRA waste code RCRA waste code		DU01 characteristic assumed due to organics. Concentrations for metals were assumed to be 200 times the regulatory level. Concentrations for organics were based on judgement.
		F005	2009	A33	/ B602	6,993	1		Senzene Pyridine Acetone Ethyl acetate n-Butanol Cyclohexenone Mathanol Xylene Ethylbenzene Ethylbenzene Ethyl ather Mathyl isobutyl ketone	50,00 10,00 10,00 10,00 10,00 10,00	0 S 0 S 0 S 0 S 0 S 0 S	CRA waste code CRA waste code colvents LDR RIA colvents LDR RIA		D001 characteristic assumed due to benzene and F005 assumed due to benzene and pyridine. All constituents included in the Solvents LDR RIA for F003 were added. Concentrations for all constituents were based on Judgement.

Appendix 6

Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RANK		RCRA weste code		1C		ource Code		orm ode	Quantity	VS	toc	Constituents	Conc. (ppm)	Constituent Source		Assumptions
73		f001 F002 F003 F005						8204	6,975	4		Irichlorosthylene letrachlorosthylene Aylene Mathylene chloride Mathylene chloride Mathyl chloroform Mathyl chloroform Island chloroform Island chlorofluoromethane Irichlorofluoromethane Acatone Chloroform Cyclohexanone n. Butanol Mathyl isobutyl ketone Phanol o Dichlorobenzene Ethyl acatate Ethylene glycol monoati Isobutanol Benzene Salenium Antimony Arsanic Barium Cadaium Chromium Copper Lead Mercury Nickel Silver Thallium Vanadium Zinc	90,000 75,000 65,000 65,000 20,000 15,000 35,000 35,000 15,000 35,000 1,000 4,500 4,500 4,500 6,05 9,000 0,05 0,05 0,05 0,05 0,05 0,05	Gen. Survey	3	Assumption used for the Generator Survey match: RCRA codes and form code.
74	1	D001 D002	,	2869	,	A37	1	8207	6,834	1		Acetaldehyde Isobutyraldehyde Methyl ethyl ketone Hydrochloric acid	5,000 5,000	BRS CAS numbers BRS CAS numbers BRS CAS numbers None	0	D001 characteristic assumed due to organics. Hydrochioric acid added to constituents to account for D002 characteristic. Concentration of HCl based on a pH of 2 to satisfy D002 characteristic. Concentration of methyl athyl ketone was obtained from the Generator Survey. Concentrations of other organic constituents were assumed to be equal to that of methyl athyl ketone. Assumption used for the Generator Survey match: exact metch on all data elements.

Appendix 6
Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RAHK		RCHA waste code	SIC Code	Source Code		form Code	Quantity	15	1,	Constituents	Conc.	(pp=)	Constituent Source	Kay	Assumptions
75			2911	A63	1	B205	6,785			1 1,2-Dichlorosthans Benzene Toluene Phenanthrene Benzo(a)anthracene Naphthalene Pyrane (hrysene Benzo(a)pyrene Othenz(a,h)anthracene Antimony Arsenic Barium Cadmium Chromium Copper Lead Mercury Nickel Selenium Silver Vanadium Zinc		1,250 1,000 100 100 100 100 70 50 100 25 20 150 30 150 40 450 450 150 2,000	RCRA waste code 1037 38 BDAT Doc. 1037-38 BDAT Doc. F037-38 BDAT Doc. F037-38 BDAT Doc. F037-38 BDAT Doc. F037-38 BDAT Doc. F037-38 BDAT Doc. F037-38 BDAT Doc. F037-38 BDAT Doc. F037-38 BDAT Doc. F037-38 BDAT Doc.	3	0001 characteristic assumed due to organics. Concentration for 1,2-dichiorocthane was based on judgement. Constituent concentrations from the F037/jtBackground Uocument, and are the midpoints of the possible concentration range for each constituent.
76	$\downarrow$	D001 D018 D025	2821	A33	1	8219	6,728	'	1	Cresols Benzene	50 50	,000 R		0	DOO! characteristic assumed due to benzene Constituent concentrations were based on judgement.
77	'	D003	2879	A32	1	B111	6,687	1	1	Hydrazina Ammonia	50 50	,000 B	RS CAS numbers RS CAS numbers	0	Judgement.  D003 characteristic assumed due to hydrazine Concentration of hydrazine obtained from the Generator Survey. Concentration of ammonia based on judgement. Assumption used for the Generator Survey match: exact match on all data elements.

Appendix 6 .

Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RANK		RCRA waste code		IC pda		ource Code		orm ode	Quantity	3.	/ fac	Constituents Conc. (ppm	Constituent Source	Key	Assumptions
78		D001 D004 D005 D006 D007 D008 D009 D010 D011 D016					4	B202	6,591	1	1	Silver   1,00	D RCRA waste code IRCRA waste code	3	Only constituents occurring in more than 33 percent of waste streams with D001 as the only R(RA code (as reported in the Generator Survey) were included to account for the D001 characteristic. For toluene and xylene the median concentrations of the wastes in the Generator Survey were used. Concentrations of metals and 2,4-D (2,4-dichlorophenoxyacetic actd) were assumed to be 200 times the regulatory levels. The waste codes and form code \$200 conflict since the waste codes do not indicate any halogenated organics while \$200 indicates the presence of halogenated organics. The RCRA codes were assumed to be more accurate than the form code.
79	,	K002		2865		A33		8203	6,554	1	,	Cumy1 pheno1	O BRS CAS numbers O BRS waste desc. O BRS waste desc. O KOO2 BDAT Doc. O KOO2 BDAT Doc. O KOO2 BDAT Doc. O KOO2 BDAT Doc.	1	Metal constituent concentrations from KOO2 BDAT background document. Concentrations for the organics were based on judgement. Concentrations for lead and chromium were adjusted since the background document indicated unrealistically high concentrations.
60	1	0001 D002 D019 D022 0027 D028 D029 D032 (003) D034		∠869		A99		8494	6,435	1	,	1,1-Dichloroethylene 1,4 Chloroform 1,2 Heachloroethane 6 Carbon tetrachloride 1 1,2-Dichloroethane 1 Hexachlorobutadiene 1 Hexachlorobutadiene 1 Hexachlorobutadiene 1 Hexachlorobutadiene 10,0	O RCRA waste code NCRA waste code O RCRA waste code O RCRA waste code O RCRA waste code O RORA waste code	2	Waste description indicated that the waste is a PCB-contaminated organic waste. 0001 characteristic assumed due to organics. Hydrochloric acid added to constituents to account for D002 characteristic. Concentration of HCl based on a pH of 2 to satisfy D002 characteristic. Concentration for PCBs was based on judgement. Concentrations for the other constituents were assumed to be 200 times the regulatory level.
01	,	F002 F003 F005	,	2384	·	A37	,	B101	6,414	1	1	Methylene chloride 1,5 Tetrachloroethylene 0	5 BRS CAS numbers 5 BRS CAS numbers 0 BRS CAS numbers 0 Gen. Survey 2 Gen. Survey	2	F003 and F005 codes assumed due to organics indicated by the BRS CAS numbers. Concentrations of these organics were obtained from the background document for F001-F005 for the fiber industry. Other constituents and concentrations obtained from the Generator Survey. Assumption used for the Generator Survey match: RCRA codes
92	,	D001 D018		2821		A33	,	8211	6,340	1		Ethylbenzene 50,0 Toluene 50,0 Xylene 50,0	BRS CAS numbers RCRA waste code	0	D001 characteristic assumed due to organics. Constituent concentrations for toluene and sylene obtained from the Generator Survey Information on D001 wastes. Other constituent concentrations based on judgement.

Appendix 6
Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RANK	RCHA weste code	SIC ode	Source Code	lorm Lode	Quantity	5.	1.4	Constituents	Conc. (ppm	Constituent Source	K	(ey	Assumptions
83	£001	2911	ASS	8603	6,217	2	2	fluorene Naphthelene Acenaphthalene Toluene Xylene Phenanthrene Anthracene Benzene fluorine Benzene fluorene Pluorenthene Pyrene Pluorenthene Pyrene Phenol 2,4-Dimethyl phenol Ethylbenzene Antimony Arsenic Barium Beryllium Cadmium Chromium Copper Lead Mercury Nickel Salenium Silver Thellium Vanadium Zinc	255 257 257 257 257 257 257 30 003 40 05 40 br>40 05 40 0	O Gen. Survey		1	Assumption used for the Generator Survey match: RIHA codes and form code.
4	f003 f005	2819	A	В	6,101	2		Methyl ethyl ketone Toluene Xylene Acetone Benzene n-Butanol Cerbon disulfide Chloroform Ethyl ether Methenol Antimony Arsenic Berfum Cadmium Chromium Lead Lead Lead Jencury Selenium Zinc	160,000 135,000 110,000 33,000 14,000 4,000 4,000 7,500 40,000 0.05 0.1	Gen. Survey	3	3 6	Assumption used for the Generator Survey match: RCRA codes and SIC code.

Appendix 6

Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RANK	Γ	RCRA weste c	ode	316 obe	ource Code		orm ode	Quantity	WS	fac	Constituents	Conc. (ppm)	Lonstituent Source	Kay	Assumptions
85	,	D001 D002 D00 D018 D021 D02 D025 D026		2869				6,075	1	1	Chlorobenzene Cresols Benzene Hydrogen sulfide Hydrochloric acid	2,000 100 500	RCRA waste code RCRA waste code RCRA waste code None None	2	DUD1 characteristic assumed from benzene and chlorobenzene and hydrogen sulfide was added to cover DUD3 characteristic. Concentrations for benzene, chlorobenzene, and cresuls are 200 times the regulatory levels. Hydrogen sulfide concentration was based on proposed EPA guidelines. Hydrochloric acid added to constituents to account for DUD2 characteristic. Concentration of HC1 based on a pH of 2 to satisfy DUD2 characteristic.
86	1	D001 F001 F00	03 F005		_	•	B204	5,956	8	8	Acetone Methanol Methyl ethyl ketone Mylane Toluene Methyl chloroform Methyl chloroform Methyl isobutyl ketone Tetrachloroethylene Carbon disulfide Chlorobanzene Chloroform Ethyl acetate Ethyl acetate Ethyl acetate Ethyl acetate I,1,2-Trichloroethane n-Butanol Isobutanol Arsenic Berium Beryllium Cadmium Chromium Copper Lead Mercury Salenium Silver Zinc	140,000 140,000 140,000 100,000 80,000 45,000 15,000 12,500 12,500 12,500 12,500 12,500 12,500 12,500 12,500 12,500 12,500 12,500 12,500 12,500 12,500 12,500 12,500 12,500 12,500 12,500 12,500 12,500 12,500 12,500 12,500	Gen. Survey	3	DOOI characteristic assumed due to organics. Assumption used for the Generator Survey match: RCRA codes and form code. Since the form code indicates a halogenated solvent mixture, concentrations were adjusted so that total concentration would be approximately 1,000,000 ppm.

Appendix 6
Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RANK	RCRA waste code	SIC Code	Seurc Code		form (ode	Quantity		141	Constituents	Conc. (ppm)	Constituent Source	Kay	Assumptions
87	1001 1001 1002 1003			,	8202	5,922	6	6	letrachlorosthylene Irichlorosthylene Hethylene chloride 1,1,1 Tichlorosthane (arbon tetrachloride (hlorobenzene 1,2 Dichlorobenzene Irichlorofluoromethane Xylene Acetone Ethyl acetate Ethyl acetate Ethyl ether Hethyl isobutyl ketone n-Butenol Cyclohexanone Hethyl ethyl ketone Isobutanol Isobutanol Pyridine	50,000 50,000 40,000 50,000 40,000 50,000 50,000 50,000 25,000 25,000 40,000 40,000 30,000 25,000	Solvents LDR RIA	2	1000) characteristic assumed due to organics. All constituents included in the LDR RIA were added. The waste code 6000 and 6000 and 6000 both indicate non-halogenated organics while 8202 indicates the presence of only halogeneted organics. The RCRA codes were assumed to be more accurate than the form code. Since the form code indicates a solvent misture, concentrations were adjusted so that total concentration would be less than 1,000,000 ppm.
68	1001 F001 F002 F003	2899	/ A89		B204	5,825	2	1	Acetone Mathylene chloride 1.1.1-Trichloroethane Mathyl chloroform Irichlorotrifluoroethane Selenium	22,000 6,000 6,000 6,000	Gen. Survey Gen. Survey Gen. Survey Gen. Survey Gen. Survey Gen. Survey	3	UIOO1 characteristic assumed due to organics. Assumption used for the Generator Survey match: RCRA codes and form code.
89	POOL DOIR	5851	A37		B219	5,724	1	1	Naphthalena Cumene Ethylbenzene Ioluene Kylene Benzene	50,000 50,000 50,000 50,000	BRS CAS numbers BRS CAS numbers BRS CAS numbers BRS CAS numbers BRS CAS numbers RCRA waste code	0	DUOI characteristic assumed due to organics. Constituent concentrations for benzene, toluene, and mylane obtained from the Generator Survey information on DUOI wastes. Other constituent concentrations based on judgement.

Appendix 6

Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RANG	Ī	RCRA	weste	code	IC ode	ource Code	ores ode	Quant ity	WS	f oc	Constituents	Conc. (ppm)	Constituent Source	Key	Assumptions .
90		0001	F003 F0	105				5,692	123	122	Toluene  Xylene  Hathyl ethyl ketone Acatone Ethyl ecetate Hethanol Hethyl chloride Hethyl isobutyl ketone Hethyl isobutyl ketone Hethyl ena chloride Acatonitrile Cyclohexanone Ethylene glycol monoeth Acatophenone Benzene Formaldehyde Isobutanol Hethyl chloroform Naphthalene Phenol Pyridine Arsenic Barium Baryllium Cadmium Chromium Copper Lead Mercury Nickel Salenium Silver Thallium Vanadium Zinc	120,000 110,000 65,000 20,000 55,000 10,000 10,000 10,000 55,500 55,500 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000 10,000	Gen. Survey	3	DOD1 characteristic assumed due to organics . Assumption used for the Generator Survey match: RCRA codes and form code.
91	,	0016		007 0008 001 F003	2869	A37	8219	5,679	1	1	Acrylonitrile Benzene Methyl ethyl ketone Chromium Lead Tatrachloroethylene Trichloroethylene Methylene chloride 1,1,2-Trichloroethane Carbon tetrachloride Kylene Acatone Ethyl acetate Ethyl setate Ethyl ether Methyl isobutyl ketone n-Butanol Cyclohexanone Methanol Hydrochloric acid	4,000 3,000 1,000 5,000 5,000 5,000 4,500 4,500 3,000 3,000 4,500 4,500 4,500	RCRA waste code RCRA waste cod		D001 characteristic and F005 code assumed due to benzene and methyl ethyl ketone. All constituents included in the Solvents LDR RIA for F001 end F003 were added. Concentrations of organics were obtained from the LDR RIA. Concentrations of metals were assumed to be 200 times the regulatory level. Concentration of acrylonitrile was based on judgement. Hydrochloric acid added to constituents to account for D002 characteristic. Concentration of HCl based on a pH of 2 to satisfy D002 characteristic. Since the form code indicates unspecified organic liquids and the BRS waste description indicated that the waste is comprised on drainage, flushings, and washings, concentrations were based on judgement and their relative proportions as given in the Solvents LDR RIA.

Appendix 6
Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RANK	Π	MCRA weste code	SIC	Source	form	Quantity	•	•	Constituents Conc. (pp	Constituent	T.	
	_		Code	Codo	Code	<u> </u>		fec	conc. (pp	Source	Key	Assumptions
92	,	D001 D004 0005 D006 D007 D008 0009 (1010 D011 D018			4 8505		5	5	Bartum   20,00	O HCRA waste code O RCRA waste code	1	D001 characteristic assumed due to benzene Constituent concentrations for metals were assumed to be 200 times the regulatory level. Benzene concentration was based on judgement. For this waste stream combination, the organic concentrations given in the background document were not used because the total organic concentration was less than 3 percent which appears to be too low for organic liquids (as is indicated by the form code). The waste codes and form code 8202 conflict since the waste codes do not indicate any halogeneted organics while 8202 indicates the presence of helogeneted organics. The RCRA codes were assumed to be more accurate than the form code.
94		(DUO1 11154	2869	A33	8219	5,554	1	1	Acrolein 25,00 Acrylamide 25,00 Acrylamide 25,00 Acrylamide 25,00 Arsenic 0.6 Barium 0.6 Nickei 0.0 Zinc 0.00	O RCRA waste code O KO13 BDAT Doc.		0001 characteristic assumed due to acrylonitrile. For this waste stream combination, the organic concentrations given in the background document were not used because the total organic concentration was less than 1 percent which appears to be too low for organic liquids (as is indicated by the form code). Instead, concentrations of organics were based on judgement with the consideration that the BRS waste description indicated that the waste is an acetonitrile waste. Concentrations for the metals were obtained from the background document.
			2869	A34	/ 8203	5,479	1	1	Methanol 560,00	RCRA waste code		DOD1 characteristic assumed due to methanol. Methanol concentration was obtained from the Generator Survey. Assumption used for the Generator Survey match: RCRA codes and form code.
95	_	D008	200			5,357	46	42	Lead 1,00	RCRA waste code		lead concentration was assumed to be 200 times the regulatory level.
		5002 0002 7003	2819	<b>A</b>	8	5,323	2		Acatona   Acatona   Acatona   Acatona   Acatona   Acatona   Antimony   O.0   Arsanic   O.0   Barium   O.0   Cadmium   O.0   Cadmium   O.0   Cadmium   O.0   Chromium   Copper   O.0   Lead   O.0   Marcus   O.0   Micha   O.0   Selenium   O.0   Silver   O.0   Thillium   O.0   Thillium   O.0   Cadmium   O.0   Cadmium	Gen. Survey Gen. Survey Gen. Survey		D001 characteristic assumed due to methanol. Hydrochloric acid added to constituents to account for D002 characteristic. Concentration of HCl based on a pH of 2 to satisfy D002 characteristic. All other concentrations were obtained from the Generator Survey. Assumption used for the Generator Survey match: RCRA codes.

Appendix 6

Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RANK	Π	RC	RA weste code		SIC Code	T	Source Code		orm ode	Quantity	vs vs		Constituents	Conc. (ppm)	Constituent Source	Key	Assumptions
97	1	DO	01 D018		264	•	A37	1	<b>B</b> 205	5,230	1	1	Benzene:	100,000	RCRA waste code	0	DODI characteristic assumed due to benzana Banzana concentration was based on judgement
98		DO	01 <b>0000</b> FU03 FOO	.5	495	3 1	A73		B20.3	4,880	3	3	lead Methyl ethyl ketone Xylene Toluene Antimony Arsenic Barium Baryllium Cedmium Chromium Copper Mercury Nickel Selenium Silver Thallium Venedium Zinc	90,000 25,000 25,000 0.003 0.003 0.03 0.003 0.3 0.003 0.3 0.	RCRA waste code Gen. Survey	1	DDD1 characteristic assumed due to organics. (uncentrations for all constituents were obtained form the Generator Survey. Assumption used for the Generator Survey match: RCRA codes and form code.
99			01 D007 D008 F00 02 F003 F005		999	9			B202	4,866	1	•	Chromium Lead Methyl chloroform Toluene Xylene Tetrachloroethylene Trichloroethylene Methylene chloride 1,1,1-Trichloroethene Carbon tetrachloride Chlorobenzene 1,2-Dichlorobenzene Trichlorofluoromethene	1,000 70,000 70,000 70,000 70,000 70,000 70,000 70,000 70,000	F001-F005 BDAT Doc. F001-F005 BDAT Doc. F001-F005 BDAT	3	D001 characteristic assumed due to organics, 5003 code assumed due to xylene, and F005 code assumed due to toluene. Concentrations for the metals were assumed to be 200 times the regulatory level. Assumption used for the Generator Survey match: RCRA codes. All organics listed under f001 and F002 were added from the F003-f005 background document. The concentrations for all organics were based on judgement. The waste code F003 and F005 and form code B202 conflict since f003 and f005 both indicate non-helogenated organics while B202 indicates the presence of only halogenated organics. The RCRA codes were assumed to be more accurate than the form code.
100		00	01 <b>D</b> 602		28:	13	A37		8101	4,830	1	1	Methanol Acatone Toluene Hydrochloric acid Sulfuric acid Isopropanol Ammonia	25,000 15,000 500 500 25,000	BRS CAS numbers BRS CAS numbers BRS CAS numbers BRS CAS numbers BRS CAS numbers BRS waste desc. BRS waste desc.	0	Organic constituents assumed to account for D001 characteristic, and acid constituents assumed to account for D002 characteristic. Concentrations of HC1 and sulfuric acid based on a pH of 2 to setisfy D002 characteristic. Concentrations for isopropanol and ammonia were based on judgement. Concentrations for other constituents were obtained from the Generator Survey. Assumption used for the Generator Survey match: RCRA codes and form code.

Appendix 6
Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RANK		RCRA weste code	S1 Cod		Source Cede		form Code	Quantity	WS	Fac	Constituents Conc. (ppm	Constituent Source	Key	Assumptions
101		F001 F002 F003						4,822	6	8	Methylene chloride   22,00    1,1,1-Trichlorosthana   6,00    Hethyl chloroform   6,00    Trichlorotrifluoroethana   6,00	Gen. Survey Gen. Survey Gen. Survey Gen. Survey Gen. Survey	3	The form code was assumed to be 8204 (mixture of halogenated and non-halogenated solvents) based on the RCRA cedes. Constituents and concentrations were obtained from waste stream combination under Rank 88
102		D001 D002 D007 D018 D021 F002 F003 F005		865	A31		B204	4,781	2	1	Mathanol   150,000   Benzene   150,000   1,2-Dichlorobenzene   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   150,000   1	BRS (AS number BRS (AS number BRS CAS number BRS CAS number RCRA waste code RCRA waste code None	3	D001 code assumed due to organics, f002 code assumed due to chlorobenzene and 1,2-dichlorobenzene, f003 code was assumed due to xylene, and f005 code assumed due to benzene. Hydrochloric acid added to constituents to account for D002 characteristic. Concentration of HCl besed on a pH of 2 to satisfy D002 characteristic. Concentration of chromium was assumed to be 200 times the regulatory level. Concentrations for all other constituents were based on judgement.
103		D001 D004 D005 D006 D007 D008 D009 D010 D016 F001		189	A71	,	8219	4,743	1	1	2,4-01chlorophenoxy acetic aci2,001	R(RA waste code R(RA waste code Solvents LDR RIA Solvents LDR RIA Solvents LDR RIA Solvents LDR RIA Solvents LDR RIA Solvents LDR RIA Solvents LDR RIA Gen. Survey	3	All constituents for FOOl were obtained from the Solvents LDR RIA. The concentrations for these organics were based on judgment since the form code unspecified organic liquids. Only constituents occurring in more than 33 percent of waste streams with DOOl as the only RCRA code (as reported in the Generator Survey) were included to account for the DOOl characteristic. For toluene and xylene the median concentrations of the wastes in the Generator Survey were used. The concentrations for the metals and 2.4 D were assumed to be 200 times the regulatory level.
104		DO01 D018 D025	29	)11	A33	ľ	8219	4,694	1	1	senzene 50.000	RCRA waste code RCRA waste code BRS waste desc.	0	D001 characteristic assumed due to benzene Concentrations were based on judgement.
105	,	K022	26	165	A35		B602	4,609	1	1	Pheno 1 50,000 Arsenic 5 Cadmium 0.5	Gen. Survey Gen. Survey Gen. Survey Gen. Survey	1	Assumption used for the Generator Survey match: RCRA codes and form code.

Appendix 6

Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RAMK		RCRA waste code	IC de	urce Cede		orm ode	Quantity	WS.	iac		Conc. (ppm)	Constituent Source	Key	Assumptions
106	i I	0001 0002 0003 10004 0005 0006 0007 0008 0009 0010	4953	A99		6)14	4,564	1	1	Barium Arsanic Laad Chromium Silver Cadmium Hercury Taluana Kylana Hydrochloric acid Hydrogen sulfide	1,000 1,000 1,000 1,000 200 40 50,000 50,000	RCRA waste code RCRA waste code RCRA waste code RCRA waste code RCRA waste code RCRA waste code RCRA waste code Gen. Survey None	1	Only constituents occurring in more than 33 percent of waste streams with D001 as the only RCRA code (as reported in the Generator Survey) were included to account for the D001 characteristic. For toluene and xylene the median concentrations of the wastes in the Generator Survey were used. The concentrations for the metals were assumed to be 200 times the regulatory level. Hydrochloric acid added to constituents to account for D002 characteristic. Concentration of HC1 based on a pH of 2 to satisfy D002 characteristic. Hydrogen sulfide included to account for D003 characteristic. Concentration of hydrogen sulfide based on proposed EPA guidelines.
107	I	D001 D004 D005 D006 D007 D008 D010 D011 D018 D035	2899	A89		B204	4,531	1	1	Hethyl ethyl ketone Barium Banzene Arsenic Lead Chromium Silver Cadmium Selenium Hercury	20,000 2,200 1,000 1,000 1,000 1,000 200	RCRA waste code RCRA waste code	1	DOO1 cheracteristic assumed due to organics. Concentrations of metals were assumed to be 200 times the regulatory lavel. Concentration of methyl athyl ketone was based on judgament.
108	,	DUO1 DO04 DO05 DO06 DO07			1	B407	4,509	4		Barium Arsenic Chromium Cadmium Toluene Xylene Lead	1,000 1,000 200 50,000 50,000	RCRA weste code RCRA weste code RCRA weste code RCRA weste code Gen. Survey Gen. Survey Gen. Survey	1	Concentrations of metals were assumed to be 200 times the regulatory level. Only constituents occurring in more than 33 percent of waste streams with D001 as the only RCRA code (as reported in the Generator Survey) were included to account for the D001 characteristic. For lead, toluene, and xylene the median concentrations of the wastes in the Generator Survey were used.
109	7	0001	2869	A33		8207	4,501	9		Formaldehyde Methanol Acetone Acetonitrile Acetaldehyde n-Butanol Ethylene glycol Ethanol	12,500 12,500 12,500 12,500 12,500 12,500	BRS CAS number BRS Weste desc.	0	DOOl characteristic assumed due to organics. Concentrations based on judgement.
110	,	K027	2865	A33	,	B403	4,457	1		Toluene-2,6-diisocyanate Toluene-2,4-diisocyanate Ethylbenzene 2,4-Toluene diamine 2,6-Toluene diamine o-Dichlorobenzene Nickel	20,000 5,000 5	BRS CAS number BRS CAS number Gen. Survey Gen. Survey Gen. Survey Gen. Survey	3	Concentrations of all constituents were obtained from the Generator Survey. Assumption used for the Generator Survey match: RCRA codes. Concentrations of the organics were adjusted to account for the form code that indicates organic solids.

Appendix 6
Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RAH		RCRA weste code	SIC Code	Saurce Code	Form Code	Quantity	vs vs	fac	Constituents .Conc. (p		Constituent Source	Key	Assumptions
111	,	0005 0006 0008 F001	4953		/ B204	4,348	1	1	Lead   1   1   1   1   1   1   1   1   1	200 000 000 000 000	RCRA waste code RCRA waste code RCRA waste code Solvents LDR RIA Solvents LDR RIA Solvents LDR RIA Solvents LDR RIA Solvents LDR RIA Solvents LDR RIA	3	All constituents (and concentrations) for f001 were obtained from the Solvents LDR HIA. The concentrations for the metals were assumed to be 200 times the regulatory level. Since the form code indicates a solvent mixture, concentrations of the organics were adjusted so that total concentration would be approximately 1,000,000 ppm
112		D001	2612	A92	/ B403	4,322	1			000 000 000 000 000 000 000 000 000 00	Gen. Survey	3	Assumption used for the Generator Survey match: RCRA codes and form code.

Appendix 6

Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RANK		RCR	A waste code	1C ode	ource Cede		orm ode	Quantity	Vs.	140		Constituents	Conc. (ppm)	Constituent Source	Kay	Assumptions
113	,	0001 0021	1 0002 (NO) 0018 1 F002 F003 F005	2865	A34		B204	4,316	1	1	8 1 0	dethanol tylene benzene i,2-Dichlorobenzene chlorobenzene chromium dydrochloric acid	100,000 100,000 100,000 2,000 1,000	BRS CAS number BRS CAS number BRS CAS number BRS CAS number RCRA waste code RCRA waste code None	3	(1001 characteristic assumed due to organics, f002 code assumed due to 1,2-dichlorobenzene, f003 code assumed due to methanol, and f005 code assumed due to benzene. Concentration of chlorobenzene and chromium was assumed to be 200 times the regulatory level. Concentrations for the organics were based on judgament. Hydrochloric acid added to constituents to account for D002 characteristic. Concentration of HCl based on a pH of 2 to satisfy D002 characteristic.
114	1	D00	1 DO18	2865	A		8	4,312	1		1 (	Benzene .	100,000	RLRA waste code	0	DOD1 characteristic assumed due to benzene. Since the waste is combusted, it is assumed to be an organic liquid. Concentration was based on judgament.
115		UOO	1 D002 D003 D035 1 U002 U008 U028 1 U057	5169	A31	•	B207	4,196	1			Methyl ethyl ketone Acataldehyde Acatone Acrylic ecid Diethylhesyl phthalata n-Butanol Cyclohexanone Hydrogen sulfide	10,000 10,000 10,000 10,000 10,000 10,000	RCRA waste code RCRA waste code RCRA waste code RCRA waste code RCRA waste code RCRA waste code RCRA waste code None	0	DOOI characteristic assumed due to organics. Concentrations of the organics were based on judgement. Hydrogen sulfide was included to account for DOO3 characteristic. Concentration of hydrogen sulfide based on proposed EPA guidelines.
116	,	100	1 D002 F002 F003 5 U002 U012 U031 4 U080	2834	A37	,	<b>6</b> 201	4,163	1			Methenol Acetone Toluene Ethylene glycol n-Butanol Methylene chloride Anlline Chloroform	100,000 100,000 10,000 100,000 10,000	BRS CAS number BRS CAS number BRS CAS number BRS CAS number RCRA waste code RCRA waste code RCRA waste code RCRA waste code	2	DOD1 characteristic, FOO2, FOO3, and FOO5 codes assumed due to organics. Constituent concentrations were based on judgement with the consideration that the BRS waste description indicated that the waste contains primarily non-halogenated solvents.
117	1	D00	1 F003	2833	A49	1	<b>B</b> 101	4,003	2			Xylene Acetone Ethyl acetate Acetonitrile Methanol	25,000 2,500 25,000	Gen. Survey Gen. Survey Gen. Survey Gen. Survey Gen. Survey	0	DOD1 essumed due to organics. Concentrations of all constituents obtained from the Generator Survey. Assumption used for the Generator Survey match: RCRA codes and form code.
118	'	000	D1 D035	2869	A35	′	8203	4,000			•	Methyl ethyl ketone	100,000	RCRA waste code	0	DOD1 characteristic assumed due to methyl ethyl hetone. Concentration based on judgement.

Appendix 6
Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

HANK	Т	RCRA weste code	Γ	SIC	Τ.	Source	T	l orm	To		Γ.	L		<u> </u>		·
	<u> </u> _		Ļ	Code		Code Code		ode	Quent 1ty	10.	140	Constituents	Conc. (ppm)	Constituent Source	Kay	Assumptions
119		6001 F004		2621		A33	1	8602	3,990	1	1	tresols 2,6-Dimethyl phenol Nitrobenzene Phenol Xylene Methylene chloride Methyl isobutyl ketone loluene Arsenic Barium Beryllium Cadmium Cropper Lead Mercury Nickel Solver Zinc	50,000 140,000 130,000 120,000 30,000 0,1 0,000 0,1 10,000 0,1 10 0,1 10 0,1 10 0,1 10 0,1 10 0,1 10 0,1	BRS (AS number BRS weste desc. Gen Survey Gen. Survey	3	ioncentration for 2,6-dimethyl phenol was based on judgement. Concentrations of all other constituents were obtained from the Generator Survey. Assumption used for the Generator Survey match: RCRA codes.
120		D001 F003		2869		A33	1	B219	3,879	2.	2	Xylene Ethylbenzene Methanol Limonene	10,000 10,000	BRS CAS number BRS CAS number BRS CAS number BRS waste desc	O	DOO! characteristic and FOO3 code assumed due to organics. Constituent concentrations were based on judgement with the consideration that the BHS weste description indicated that the waste is a mixture of terpenes containing small amounts of sylene.
122	·	D001 D002 D019 D032		2869		A31		8207	3,873	1	1	Mathanol Copper Hexachlorocyclopentadie Hydrochloric acid	15,000 ne 5,000	Gen. Survey Gen. Survey Gen. Survey None	j	Assumption used for the Generator Survey match: exact metch on all data elements. Hydrochloric acid added to constituents to account for DOO2 characteristic. Concentration of HCl based on a pH of 2 to satisfy DOO2 characteristic. Concentration for copper was adjusted since the Generator Survey indicated an unrealistically high concentration for copper.
123		D033 D034 D039 F002		2869		A33		8219	3,866	2	1	Carbon tetrachloride Hexachlorobenzene Hexachlorobethene Tetrachloroethene Tetrachloroethylene Chloropyridine Toluene Xylene Lead Chromium Cadmium Hydrochloric acid	50,000 50,000 50,000 50,000 150,000 50,000 50,000 50,000	RCRA waste code RCRA waste code RCRA waste code RCRA waste code BRS waste desc. Gen. Survey Gen. Survey Gen. Survey Gen. Survey Gen. Survey None		F002 code assumed due to tetrachlorosthylane. Only constituents occurring in more than 33 percent of weste streams with D001 as the only RCRA code (as reported in the Generator Survey) were included to account for the D001 characteristic. For the metals the median concentrations of all the D001 westes in the Generator Survey were used. The concentrations of all organics were besed on judgement. Hydrochloric acid added to constituents to account for D002 characteristic. Concentration of HCl based on a pH of 2 to satisfy D002 characteristic.
123		D001		2865		A33	1	<b>B</b> 203	3,852	1	1	Acetone Benzene Cumene Phenol e Methyl styrene Ethylbenzene	80,000 80,000 80,000 80,000	BRS CAS number BRS CAS number BRS CAS number BRS CAS number BRS waste desc. BRS waste desc.	0	DOO1 characteristic assumed due to organics. Constituent concentrations were based on judgement.

Appendix 6

Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

					 			<del></del>	_			<del></del> - 1		1	Assumptions
RANK		RCRA weste code	S1 Co		ource Code		or <b>u</b>	Quantity	WS	fac	Constituents Con	c. (ppm)	Constituent Source	,	ASSEMPTIONS
124	1	DOO1 DOO5 DOO6 DOO?					B204	3,775	1	1	Barium Cadmium Chromium Ioluene Kylene Lead	200 1,000 50,000 50,000	RCRA waste code RCRA waste code RCRA waste code Gen. Survey Gen. Survey Gen. Survey	1	Concentrations for metals indicated by the RCRA codes were assumed to be 200 times the regulatory level. Only constituents occurring in more than 33 percent of weste streams with D001 as the only RCRA code (as reported in the Generator Survey) were included to account for the D001 characteristic. For lead, toluene, and sylene the median concentrations of the wastes in the Generator Survey were used.
125		D002 D006		2833	A32	•	8207	3,724	1	1	Cadmium Selanium Hydrochloric acid	1	RCRA waste code Gen. Survey None	1	Concentrations for the metals were obtained from the Generator Survey. Assumption used for the Generator Survey match: RCRA codes and source code. Hydrochloric acid added to constituents to account for D002 characteristic. Concentration of HCl based on a pH of 2 to satisfy D002 characteristic. Though the form code indicates an organic liquid, no organics were added with the assumption that in this case the RCRA codes were more accurate than the form code.
126	,	D001 D018 K048 K049		2911		,	B204	3,669	1	1	Benzene Lead Chromium		RCRA waste code KO48-51 List. Doc. KO48-51 List. Doc.	1	Benzene assumed to account for all RCRA codes; lead and chromium were added based upon the K048-51 listing document. Concentrations of the metals were assumed to be 200 times the regulatory level. Benzene concentration was based on judgement. Assumed that no helogenated organic constituents were present in this weste although the form code indicates the potential presence of helogenated organics.
12!	+	0018	╂╌┨	2869	A33	7	8219	3,599		1	Benzene	100,000	RCRA weste code	0	Concentration based on judgement.
120	17	0001	П	2869	A89	1	B205	3,563	1		Hethyl methacrylate	50,000	Gen. Survey	l °	Assumption used for the Generator Survey match: RCRA codes and form code.
129		D001 D005 D006 D007 D008 F001 F002 F003 F004 F005		7389	A71	1	B206	3,518	1		Barium Cadmium Chromium Lead Xylene Toluene 1,1,1-Trichloroethane Benzene Tetrachloroethylene Trichlorotrifluoroethane Haphthalene Trichlorodifluoromethane Benzo(a)anthracene Benzo(a)pyrene Cresois Nitrobenzene	5 30 90 3,300 3,100 1,200 1,100 800 600 500 250 250	RCRA waste code RCRA waste code RCRA waste code RCRA waste code Used Oil RIA Used Oil RIA FOOL-FOOS BDAT Doc: FOOL-FOOS BDAT	3	DOOI characteristic assumed due to organics. Crasols and nitrobenzene were added to account for the FOO4 code. The concentrations for these two organics were based on judgement. All the other organic constituents and concentrations were obtained from Table V-38, "Mean Concentrations of Potentially Hazardous Constituents in Used Oil Burned as Onspecification fuel," in the used oil RIA These constituents were assumed to account for FOOI-FOO3 and FOO5 codes. Concentrations for metals were obtained from the used oil RIA.

Appendix 6
Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RANK	T	RCRA wasta code	SIC	T	Source		0170	Quantity	•	,	Constituents	Conc. (ppm)	I	r	
130	+-	DU01 D002	Code	┿	Code	-	ode		ws	Fac		CONC. (PPE)	Constituent Source	Key	Assumptions
131	,	D001 F003 F005	305.3					3,477	12	10	Toluene Kylene Lead Chromium Cadmium Hydrochloric acid	50,000 5 5 0.1	Gen. Survey Gen Survey Gen Survey Gen. Survey Gen. Survey None		Only constituents occurring in more than Jipercent of waste streams with DOOI as the unity RIRA code (as reported in the sensestor Survey) were included to account for the DOOI characteristic. For these constituents the median concentrations of the wastes in the Generator Survey were used. Hydrochloric acid added to constituents to account for DOO2 characteristic. Concentration of HCl based on a pH of 2 to satisfy DOO2 (characteristic.
132		(1001 too22	2869		A37		8403	3,465	1	1	Kylene Toluena Hethyl ethyl ketone Methyl isobutyl ketone Arsenic Barium Cadmium Chromium Copper Lead Hercury Nickel Selenium Silver Zinc	70,000 10,000 10,000 10,000 11,000 0.1 0.1 125 0.01 0.1 0.1	Gen Survey Gen, Survey	1	Assumption used for the Generator Survey match: RCRA codes and form code.
					A3/		B202	3,414	1	-	Chloroform Lead Chromium Cadmium	5 5	BRS CAS number Gen. Survey Gen. Survey Gen. Survey		Only constituents occurring in more than 33 percent of weste streams with D001 as the only RCRA code (as reported in the Generator Survey) were included to account for the D001 characteristic. For the metals the median concentrations of the wastes in the Generator Survey were used. Since the form code indicates a halogenated solvent mixture, the non-halogenated organics included in the Generator Survey information on D001 wastes were not added. Chloroform concentration was based on
133	1	D001 D005 D006 D007 D008 F003 F005	2821	,	A73	1	0602	3,410	1		Barium Cadmium Chromium Lead Toluene Xylene	1,000 1,000 50,000	RCRA waste code RCRA waste code RCRA waste code RCRA waste code Gen. Survey Gen. Survey	1	Judgement.  D001 code assumed due to mylene, F003 code assumed due to mylene, and F005 code assumed due to toluene. Concentrations for metals were assumed to be 200 times the regulatory level. Concentrations of the organics were obtained from the Generator Survey. Assumption used for the Generator Survey match: RCRA codes.

Appendix 6

Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RANK	Γ	RCPA waste code		IC do	ource Code	ore ode	Quantity	<b>%</b>		Constituents Conc. (ppm)		Constituent Source	Key	Assumptions
134	,	KO48 KO49 KO51					3,393	3		Senzo(1) fluorenthene	000000000000000000000000000000000000000	Gen. Survey	1	Assumption used for the Generator Survey match: RCHA codes.
135		E049	1	2911	A75	8202	3,316	1		Bis (2-ethylhesyl) phthelate   700	0000500000002511010102110	Gen. Survey	3	Assumption used for the Generator Survey match: RCRA codes, SIC code, and source code.

Appendix 6

Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RANK		RCRA weste code		SIC Code	I	Source Code		form Code	Quantity	vs.	, ac	Constituents	Cosc. (ppm)	Constituent Source	Key	Assumptions
136	,	D001 D002 D005 D006		495.			1	<b>8</b> 204	3,295	1	1	Barium Cadmium Toluene Xylana Lead Chromium Hydrochloric acid	200 50,000 50,000 5	RCRA waste code RCRA waste code Gen. Survey Gen. Survey Gen. Survey None	1	(oncentrations for metals indicated by the 10 codes were assumed to 200 times the regulatory level. Only constituents occurring in more than 33 percent of waste streams with DOD1 as the only RCRA code (as reported in the Generator Survey) were included to account for the DOD1 characteristic. For these constituents the median concentrations of the wastes in the Generator Survey were used. Hydrochloric acid added to constituents to account for DOD2 characteristic. Concentration of HC1 based on a pH of 2 to satisfy DOD2 characteristic. Assumed that no helogenated organic constituents were present in this waste although the form code indicates the potential presence of helogenated organics.
1 16	•	0001 D008 D008 1002	,	2869		A35	1	8403	3,175	2	1	Benzene Chlorobenzene o-Dichlorobenzene p-Dichlorobenzene	5,000 300	BRS waste desc. Gen. Survey Gen. Survey Gen. Survey	2	Benzene was added since the BRS waste description indicated that the waste is burned for energy recovery and the Generator Survey indicated the presence of chlorobenzenes. Concentration for benzene was based on judgement. Concentrations for the other constituents were obtained from the Generator Survey. Assumption used for the Generator Survey match: exect match on all data elements.
									3,,,68	۷.	2	Cadmium Lead Tolusne Xylene Chlorobenzene Trichlorofluoromethane Irichloroethylene Tetrachloroethylene Hethylene chloride 1,1,1-Trichloroethane 1,2-Dichlorobenzene	1,000 1,000 1,000 1,000 1,000 1,000 1,000	ACRA waste code RCRA waste code Gen. Survey Gen. Survey Solvents LDR RIA Solvents LDR RIA	3	Concentrations for metals were assumed to 200 times the regulatory level. Only constituents occurring in more than 33 percent of waste streems with D001 as the only RCRA code (as reported in the Generator Survey) were included to account for the D001 characteristic. For cadmium and lead the median concentrations of the wastes in the Generator Survey were used. Constituents included in the Solvents LDR RIA for F002 were added. Concentrations
39		D019 D022 D032 D039 D043 K018 K020		2869		A33	,	8219	3,132	1.		Chloroform Tetrachloroethylene Carbon tetrachloride Vinyl chloride Hexachlorobenzene Ethylene dichloride	20,000 20,000 20,000 20,000	RCRA waste code RCRA waste code RCRA waste code RCRA waste code RCRA waste code BRS waste desc.	i	for the organics were based on judgement.  KO18 code assumed due to hexachlorobenzene and KO20 code assumed due to tetrachloroethylene. Concentrations for all constituents were based on judgement
40	- 1	D001 D005 D006 D007 D008 D011 D022 D035 D039 F001				-			3,124	1		Methyl ethyl ketone Barium Chromium Lead Silver Chloroform Cadmium Tetrachloroethylene	40,000 20,000 1,000 1,000 1,000 1,000 200	RCRA wasta code RCRA wasta code		DOOI characteristic assumed due to methyl estimate and FOOI code assumed due to tetrachloroathylane. Concentrations for all constituents were assumed to be 200 times the regulatory level.

Appendix 6

Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RANK	T	R(RA waste code	SIC ode	ource Cede		or <b>m</b> ode	Quantity	VS	fac	Constituents	Conc. (ppm)	Constituent Source	Kay	Assumptions
141	1	0001 F003 F005	2869	A09	1	8203	3,071	,	3	Toluene Methanol Ethanol	250,000	BRS CAS number BRS CAS number BRS waste desc.	0	FOO3 code assumed due to methanol and FOO5 code assumed due to toluene. Constituent concentrations were based on judgement.
142	,	0001	2822				3,068	ì		Styrene Toluene Hethyl athyl ketone Carbon tetrachloride 1,1-Dichlorosthylene Kylene Acrylonitrile Chloroprene Acrylonitrile Chloroprene Ethyl acetate Ethyl acetate Ethyl acetate Ethyl henzene Hethenol Hethyl chloride Hethyl chloride Banzene Antimony Arsenic Barium Beryllium Cadmium Copper Lead Mercury Nickel Salenium Silver Thallium Vanadium Zinc	15,000 5,000 4,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 2,500 1,000 1,000 2,500 1,000 1,000 2,500 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000 1,000	Gen. Survey	3	Assumption used for the Generator Survey match: RCHA codes and SIC code.
143	1	0019 D038	2869	A33	′	B606	3,060	1	1	Benzene Pyridine		RCRA waste code RCRA waste code	0	Constituent concentrations based on judgement.
144		D001 D005 D006 D007 D008 D009 D019 F001 FQ02 F003				B407	2,999	1	1	Barium Chromium Lead Cadmium Carbon tetrachlorida Marcury Tetrachloroethylene Irichloroethylene Hathylene chloride 1,1,1-Trichloroethane Carbon tetrachlorida Chlorobenzene 1,2-Dichlorobenzene 1,2-Dichlorobenzene Erichlorofluoromethane Xylene Acatone Ethyl acatate Ethyl acatate Ethyl sther Methyl isobutyl ketone n-Butanol Cyclohexanone Mathanol	1,000 200 100. 40 3,000 3,000 3,000 3,000 3,000 3,000 3,000 3,000 3,000 3,000 3,000	RCRA waste code RCRA waste code RCRA waste code		DOOI assumed due to organics and fOOI assumed due to carbon tetrachloride. Constituents given in the Solvents LDR RIA for FOO2 and FOO3 were added. Concentrations for metals were assumed to 200 times the regulatory level. Concentrations for other constituents were hased on judgement.

Appendix 6
Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RANK		RCRA weste code	IC de		urce ode		orm ode	Quantity		100	Constituents	Conc. (ppm)	Constituent Source	Kay	Assumptions
145	,	UUU1 FAO1 FOO5	2869	^	133	,	B204	2,976	5	┾═	Toluene 2-Chloro-1,3-butadiene Tetrachloroethylene Irichloroethylene Hethylene chloride 1,1,1-Trichloroethene Carbon tetrachloride	100,000 100,000 100,000 100,000 100,000	BRS CAS numbers BRS CAS numbers Solvents LDR RIA Solvents LDR RIA Solvents LDR RIA Solvents LDR RIA Solvents LDR RIA	2	DODI characteristic assumed due to organics and 1005 code assumed due to toluene. All constituents included in the Solvents LDR RIA for FOOI were added. Concentrations for all constituents were based on judgement.
146	1	D001 D002	9999	٨			В	2,969	2	1	Toluene Xylene Lead Chromium Cadmium Hydrochloric acid	50,000 50,000 5 5	Gen. Survey Gen. Survey Gen. Survey Gen. Survey Gen. Survey None	1	Only constituents occurring in more than 33 percent of waste streams with DOO1 as the only RCRA code (as reported in the Generator Survey) were included to account for the DOO1 characteristic. For these constituents the median concentrations of the wastes in the Generator Survey were used Hydrochloric acid added to constituents to account for DOO2 characteristic. Concentration of HC1 based on a pH of 2 to satisfy DOO2 characteristic.
148		D001 D002 D003 D007	2869		37	-	8219	2,947	1		Methanol Isoheptene	500,000 100,000	BRS (AS numbers BRS waste desc.		0001 characteristic assumed due to methanol. Concentration for methanol obtained from the Generator Survey. Assumption used for the Generator Survey match: RCRA codes and form code. Concentration for isoheptane was based on judgement.
		1018 D019 (1021 D022 D028		A	3.3	′	B212	2,900	1		1,2-Dichloroethene Chlorobenzene Benzene Carbon tetrachloride Chloroform Chromium Hydrochloric acid Hydrogen sulfide	20,000 20,000 20,000 20,000 1,000 500	BRS CAS numbers BRS CAS numbers RCRA waste code RCRA waste code RCRA waste code RCRA waste code RCRA waste code None	-	D001 characteristic assumed due to benzene Concentration of chromium was assumed to be 200 times the regulatory level. Concentrations of the organics were based on judgement. Hydrochloric acid added to constituents to account for D002 characteristic. Concentration of HCl based on a pH of 2 to satisfy D002 characteristic. Hydrogen sulfide included to account for D003 characteristic. Concentration of hydrogen sulfide based on proposed EPA guidelines.

Appendix 6

Top 150 Routinely Generated Combusted Wastes: Constituent and Concentration Listing (continued)

RANK	T	RCRA weste code	)[ ode	ource Code		or <b>m</b> ode	Quantity		100	Constituents Co	nc. (ppm)	Constituent Source	Key	Assumpt fons
149		(2001 (2002	2821	A49	1	6102	2,868		,	Ethylene glycol Toluene Xylene Lead Chromium Cadmium Ethyl acetate Hydrochloric acid	25,000 25,000 5 0 1 25,000	BRS CAS numbers Gen. Survey Gen. Survey Gen. Survey Gen. Survey Gen. Survey None None	1	Unly constituents occurring in more than 13 percent of waste streams with DOO1 as the unity RINA code (as reported in the Generator Survey) were included to account for the BUO1 cheracteristic. For these constituents the median concentrations of the wastes in the Generator Survey were used. Hydrochloric acid added to constituents to account for DOO2 characteristic. Concentration of HC1 based on a pil of 2 to satisfy DOO2 characteristic. Ethyl acatate was added since the BRS waste description indicated that the waste was generated from the manufacture of polyester resins. Concentrations for ethyl acatate and athylene glycol were based on judgement Since the form code indicates an aqueous liquid, concentrations were adjusted so that total concentration would be approximately 100,000 ppm.
150	1	D002	2821	A 33	1	8105	2,832	1	,	Hydrochloric acid	500	None .	o	Hydrochloric acid was added to account for 1002 characteristic. Concentration of HCl based on a pH of 2 to satisfy D002 characteristic.

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# APPENDIX 7 SUMMARY OF PRIORITIZATION SYSTEMS

# Arizona Waste Minimization Project: Analysis of the Facility Annual Reports from the 53 Largest Hazardous Waste Generators in the State of Arizona (Task 2)<sup>1</sup>

## Purpose of the Methodology/System

- Purpose of Arizona Waste Minimization Project is to allocate state's resources towards wastes and industries with the greatest potential to succeed at further waste minimization efforts.
- Purpose of Task 2 of the Project is to identify industries that generate the largest waste volumes, and wastes that are the most significant in terms of volume and toxicity.

#### System Description

- Grouped hazardous wastes from Arizona's 53 largest waste generators into 13 "waste categories" based on generating process, chemical composition, EPA hazardous waste number, physical form, and source codes.
- Assigned relative "toxicity" scores to waste categories reflecting "threat to ground-water resources" using a ranking system based primarily on professional judgment:
  - "1" = constituents would not migrate easily to ground water
  - "2" = corrosive wastes with high levels of soluble metals
  - "3" = wastes composed primarily of solvents
- Waste categories with large volumes and high relative toxicity scores were identified as being of highest priority for waste minimization (e.g., waste categories with volumes > 100,000 pounds and a toxicity score = 3 are classified as "high priority" for waste minimization).
- Report also examines proportion of waste already recycled for each industry and waste, and reduction in waste quantity between 1989 and 1990; information will be used in later tasks to identify industries and wastes that will benefit most from further waste minimization efforts.

#### System Evaluation

Arizona Waste Minimization Project: Analysis of the Facility Annual Reports from the 53 Largest Hazardous Waste Generators in the State of Arizona, August 1992, Report prepared for U.S. EPA Office of Waste Programs Enforcement by PRC Environmental Management, Inc.

Contact: William Wilson (415/744-2153)

<sup>&</sup>lt;sup>1</sup> Documentation reviewed:

## Scope of coverage

- Hazardous wastes generated by the 53 largest generators in Arizona, excluding:
  - Wastes generated by one-time incidents
  - Wastes coded as being generated by commercial TSDFs.

## Media/pathways addressed:

Ground water

## Types of targeting criteria used (see Exhibit 1):

- Risk based:
  - Quantity of waste generated
  - Judged "threat to ground-water resources"
- Non-risk based:
  - Later tasks: potential to succeed at further waste minimization efforts (report indicates that this is related to proportion of waste already recycled)

#### Data requirements:

- Voiume of waste generated
- Waste characteristics (chemical composition, physical form, source and generating process)
- Waste management practices (proportion currently recycled)
- Data derived from Facility Annual Reports (FARs) submitted to the Arizona Department of Environmental Quality by large quantity hazardous waste generators in Arizona.

# Applicability to waste minimization targeting:

#### Pros

- Based on a small number of targeting criteria.
- Consideration (in subsequent tasks) of quantity of waste already recycled will avoid targeting industries that are already practicing effective waste minimization.

#### Cons

- Grouping wastes into categories and assigning toxicity scores relies on professional judgment; methods appear rudimentary and are not well documented. Furthermore, the 'toxicity' score does not appear to be an accurate measure of the constituents inherent toxicity.
- At the national level, availability of data for grouping wastes and assigning toxicity scores will be limited.
- Addresses only ground water.

Exhibit 1. Targeting Criteria Used in the Arizona Waste Minimization Project

TARGETING CRITERIA	
Direct Risks	
Waste volume	1
Waste/constituent toxicity Human toxicity Ecological toxicity	
Constituent concentration or mass	
Waste type (e.g., solvents)	1
Number of generators	
Waste management practices (i.e., to determine if recycling is in practice)	1
Releases to environmental media (only potential for release)	1
Potential for constituent transport	1
Potential for cross-media transfer	
Potential for human exposure: residential settings occupational settings	
Potential for ecological exposure	
Indirect or Acute Risks	
Ignitability, corrosivity, reactivity	
Ozone depletion potential	
Global warming	
Other Targeting Criteria	
Hazardous waste management capacity	
Technical/administrative feasibility	
Permitting/enforcement factors	
Cost savings	
Other factors Proportion of waste already recycled	1

# EPA/OPPT Chemical Use Clusters Scoring Methodology<sup>2</sup>

## Purpose of the Methodology/System

- To provide a screening-level scoring/ranking of chemicals identified for a specific use cluster. (A chemical use cluster refers to a group of chemicals that are all associated with a specific process or function, be it chemical, industrial, or agricultural [e.g., vulcanizing agents in the rubber chemical production process, or pesticides for a specific crop].)
- To identify use clusters that may have high potential for pollution prevention.
- Applications identified in the documentation include:
  - "to systematically identify and screen concerns related to a greater number of chemicals in commerce;" and
  - "to provide an initial indication of potentially safer substitutes for extremely toxic chemicals."

## System Description

- Two distinct components:
  - the first component scores chemicals within a single use cluster using five criteria
  - the second component scores the entire use cluster based on the chemical-specific scores and one additional criterion.
- Each of these six criteria includes one or more subcriteria which are scored based on chemical-specific data to form the basis of the numerical scores assigned to the criteria; the criteria and subcriteria are listed in Exhibit 1.
- Four scoring steps lead to the derivation of the overall chemical score:
  - based on chemical-specific data, each subcriterion is scored as high-, medium-, or low-concern, which translate into numeric scores of 3, 2, or 1.
  - the single highest subcriterion score is assigned to the criterion (e.g., human hazard potential) being scored. To minimize or "dampen" the effect of missing data, all subcriteria are regarded as equally important, i.e., equal weight, and the highest

Chemical Use Chisters Scoring Methodology, April 13, 1993 Draft Report, prepared by the Chemical Engineering Branch, Office of Pollution Prevention and Toxics, USEPA.

Contact: Daniel Fort (202/260-1694)

<sup>&</sup>lt;sup>2</sup> Documentation reviewed:

Exhibit 1: Criteria and Subcriteria

	Criteria	Subcriteria
Chemical-level	Human exposure potential	Chemical use volume Total TRI release Consumer use No. of workers No. of use sites Bioaccumulation Persistence
	Human hazard potential	Noncancer effects Cancer effects
	Ecological exposure potential	Chemical use volume Total TRI release Consumer use No. of use sites Bioaccumulation Persistence
	Ecological hazard potential	Aquatic toxicity
	EPA regulatory interest	(No subcriteria)
Cluster-level	Pollution prevention potential	Ecological risk reduction potential Human health risk reduction potential Chemical release reduction potential

score produced by any one subcriterion is assigned as the criterion score for the chemical.

- the scores of the two human risk-related criteria, i.e., human hazard potential and human exposure potential, are multiplied to obtain a score for a "secondary" criterion called the chemical human risk reduction potential.
- similarly, a score is obtained for another secondary criterion, the chemical ecological risk reduction potential.
- the overall chemical score is calculated as the sum of a chemical's scores for the following criteria:
  - human risk reduction potential;
  - ecological risk reduction potential; and
  - EPA regulatory interest.

## Derivation of the final cluster score:

- -- Score based on:
  - the individual member chemicals' overall scores
  - the score for one cluster-level criterion, i.e., the cluster's pollution prevention potential.
- pollution prevention potential criterion is scored based on three subcriteria:
  - the cluster human risk reduction potential;
  - cluster ecological risk reduction potential; and
  - chemical release reduction potential.
- the final cluster score is derived as the sum of the pollution prevention potential score and the mean of all overall chemical scores for that cluster.

## System Evaluation

## Scope of coverage:

Addresses individual TSCA chemicals within the context of their use cluster. (That is, data for some of the criteria are dependent on the specific use cluster being examined, and may differ for a given chemical with several uses.)

#### Media addressed:

- System does not score the chemical on a media- or pathway-specific basis; only aggregate releases to the environment are considered.
- However, because it uses total release information based on the Toxics Release Inventory (TRI), this system implicitly addresses the following media:
  - ... direct releases to air
  - land (landfill, underground injection)
  - surface water (direct releases, POTW transfers)

Cross-media transfers are not explicitly considered.

# Types of targeting criteria used:

- Criteria and criteria-score combinations used in this system generally conform with the risk assessment paradigm (i.e., hazard x exposure = risk).
- Of the chemical-level criteria listed in Exhibit 1, two relate to human risk potential (i.e., human hazard potential and human exposure potential), and another two to ecological risk potential (i.e., ecological hazard potential and ecological exposure potential):

- "EPA's regulatory interest" criterion assigned as follows:
  - chemical is directly assigned a score of 3, 2, or 1 for this criterion based on the number of regulatory lists of interest on which it appears (e.g., CWA Priority Pollutants list; RCRA P and U lists).
  - criterion meant to measure the Agency's previous and current interest in the chemical.
- "Cluster's release reduction potential (RRP)" criterion (the cluster's RRP score is the average of the RRP scores for all chemicals in that cluster):
  - based on its member chemicals' "efficiency-of-use," i.e., ratios of their release to use volume, estimated from TRI data (this ratio measures the fraction of a chemical throughput that is released to the environment).
- Key targeting criteria are summarized in Exhibit 2.

## Data requirements:

- Type of data required varies considerably across criteria/subcriteria, but data requirements
  are moderate.
- Use of specified data sources:
  - reviewed documentation specifies a data hierarchy for most of the criteria/subcriteria; sources of input data are well described (e.g., TRI, Clinical Toxicology of Commercial Products, HEAST, IRIS)
- Specifies methods for handling missing data and the use of professional judgment.

# Applicability to waste minimization targeting:

#### Pros

- System explicitly considers pollution prevention potential of chemicals.
- Although data-intensive, system is flexible in using data of varying quality, and allows use
  of professional judgment

#### Cons

- Focuses on the "riskiness" of a chemical specific to a use.
- System uses a number of exposure-related subcriteria that are not directly comparable. Some of the exposure-related subcriteria measure exposure magnitude (e.g., total

releases), others exposure extent (e.g., number of workers), and still others chemical properties related to exposure likelihood, magnitude, and duration. This part of the system may be internally inconsistent (since exposure is assigned a score using the single highest-scoring subcriterion) and may also "double-count" exposure, and as a result, exposure may be weighted too high relative to toxicity. That is, because there are similar exposure factors in both the human ecological criteria, exposure is essentially "double counted" relative to hazard.

Exhibit 2. Targeting Criteria Used in the Chemical Use Clusters Scoring Methodology

TARGETING CRITERIA	
Direct Risks	
Waste volume	
Waste/constituent toxicity: Human toxicity Ecological toxicity	1
Constituent concentration or mass	1
Waste type (e.g., solvents)	
Number of generators (use sites)	1
Waste management practices	
Releases to environmental media	. 1
Potential for constituent transport	1
Potential for cross-media transfer	
Potential for human exposure: residential settings occupational settings	11
Potential for ecological exposure	1
Indirect or Acute Risks	
Ignitability, corrosivity, reactivity	
Ozone depletion potential	
Global warming	
Other Targeting Criteria	
Hazardous waste management capacity	
Technical/administrative feasibility	
Permitting/enforcement factors	
Other factors:  EPA regulatory interest pollution prevention potential	1

# EPA's 33/50 Program<sup>3</sup>

## Purposes of the Methodology/System

- To target 17 chemicals and reduce their national aggregate releases by 33% by the end of 1992 and 50% by the end of 1995.
- To encourage pollution prevention activities, including source reduction and in-process recycling, in achieving these reductions.
- A voluntary reduction program. Not legally enforceable; companies are free to commit to their own reduction goals and to develop their own cost-effective strategies.
- EPA will not measure individual company efforts in program, but instead will measure progress on a national, aggregate basis (i.e., looking at the reduction in total releases of all 17 chemicals).

#### System Description

- In the targeting process, the 33/50 Program:
  - canvassed all major EPA offices for a list of their highest priority chemicals based on each office's own ranking criteria and selected from the TRI (a public data base containing information on annual releases and transfers of about 300 toxic chemicals);
  - designated all chemicals in more than one list as potential candidates for the program (25 chemicals); and
  - narrowed down the 25 chemicals to 17 by informally applying three criteria:
    - production and environmental release volumes;
    - toxicity to humans; and
    - the potential for reducing releases through pollution prevention practices.

The 17 chemicals were selected from a pool of TRI chemicals; the TRI will be used to track the progress of reaching the 50% reduction goal.

U.S. EPA. 1992. Questions and Answers: U.S. EPA's 33/50 Program [no document number given].

Telephone conversations with EPA staff members of the 33/50 Program: David Sarokin (Project Manager), John Harman, Mike Burns, Loren Hall.

Contact: David Sarokin (202/260-6396)

<sup>&</sup>lt;sup>3</sup>Documentation reviewed:

- Process used in targeting chemicals was qualitative and did not rely on a standard targeting approach:
  - -- 17 chemicals are not necessarily the "riskiest"
  - -- 17 chemicals are not the only chemicals the program is concerned about
  - -- the program, if successful, will be expanded to cover additional chemicals
- The 'narrowing down' process had the following characteristics:
  - -- no numerical ranking; ranked toxicity and volume in a high-medium-low system
  - -- professional judgment used
  - no standardized measure used for volume, though total releases and transfers in TRI were commonly used by EPA offices
  - -- no standardized measure used for toxicity, though RQs and carcinogenic potential were commonly used by EPA offices
  - pollution prevention potential criterion was especially subject to professional judgment, based on "the collective experience of the people" involved with the program
  - most weight was given to pollution prevention potential because most of candidates were high-volume and high-toxicity chemicals

#### System Evaluation

# Scope of coverage:

All constituents found in the TRI.

## Media addressed:

- The 17 target chemicals were selected partially because they are produced in large quantities and released in large quantities into the various media addressed in the TRI, including:
  - mir releases
  - surface water releases
  - discharge to POTWs
  - on-site land releases
  - off-site transfers

## Types of targeting criteria used:

• Quantity of production and environmental release (total TRI releases and transfers)

- Toxicity/hazard (RQs and carcinogenic potential)
- Potential for reduction through pollution prevention practices
- Individual offices used own targeting criteria that are not specified in the documentation reviewed
- Targeting criteria used in the 33/50 Program are summarized in Exhibit 1.

# Applicability to waste minimization targeting:

#### Pros

- Targeting process incorporated perspective from major EPA offices
- The 33/50 Program's concept of using a 'pilot' phase, i.e., targeting a few high-risk (but not necessarily the highest risk) candidates and then expanding the program if successful, may be applicable to developing the waste minimization targeting program.

#### Cons

• The program's targeting process was heavily dependent upon professional judgment and was not scientifically rigorous.

Exhibit 1. Targeting Criteria Used in the 33/50 Program.

TARGETING CRITERIA	
Direct Risks	
Waste volume	
Waste/constituent toxicity Human toxicity Ecological toxicity	,
Constituent concentration or mass	<i>J</i> 1
Waste type (e.g., solvents)	
Number of generators	
Waste management practices	
Releases to environmental media	j
Potential for constituent transport	
Potential for cross-media transfer	
Potential for human exposure: residential settings occupational settings	
Potential for ecological exposure	
Indirect or Acute Risks	
Ignitability, corrosivity, reactivity	
Ozone depletion potential	
Global warming	·
Other Targeting Criteria	
Hazardous waste management capacity	
Technical/administrative feasibility	
Permitting/enforcement factors	
Cost savings	
Other factors Pollution prevention potential Canvass of EPA offices	<i>J</i> <sub>2</sub>

Mass released to various media.
 System used chemicals determined as being of priority by other EPA offices.

# EPA Regional Comparative Risk Ranking Program<sup>4</sup>

#### Purpose of the methodology/system

- Analytical framework to systematically measure, compare, and rank environmental problems to:
  - -- help risk managers identify the worst environmental problems and risks;
  - -- provide common ground for evaluating the net benefits and costs of different strategies for reducing or preventing risks; and
  - share acquired information with the community and general public.

#### System description

- Although the system provides fairly specific guidance and methods, it allows flexibility in implementing the approach.
- System has three main components: (1) project planning and start-up; (2) comparative risk ranking; and (3) risk management.
- Comparative risk ranking is divided into four parts: separate rankings of risks to human health, ecosystems, and quality of life; and an integrated ranking of these three risk areas. The integrated ranking groups the problems in several tiers of descending risk.
- The system components are summarized in Exhibit 1. The component for assessing risks to human health, ecosystems, and quality of life is described below.
  - (1) Assessing risks to human health
  - Identify hazards to human health, including:
    - target pollutants (limit to pollutants that best represent actual risks)
    - -- relevant exposure pathways
    - adverse health effects (cancer and/or non-cancer)

U.S. EPA. September 1993. A Guidebook to Comparing Risks and Setting Environmental Priorities. Prepared by the Office of Policy, Planning, and Evaluation, U.S. Environmental Protection Agency, Washington, D.C. EPA 230-B-93-003.

Vermont Agency of Natural Resources. Environment 1991: Risks to Vermont and Vermonters. Report by the Public Advisory Committee, The Strategy for Vermont's Third Century.

Contact: Debora Martin (202/260-2699)

<sup>&</sup>lt;sup>4</sup> Documentation reviewed:

# Exhibit 1: Components of the Comparative Risk Ranking Program

#### (1) Project Plausing and Start-up

- Assemble planning team:
  - Project Manager
  - Steering Committee
  - Public Advisory Committee
  - Technical Work Groups
- Define scope of and goals for the project
- Secure support of key stakeholders
- Determine public participation role
- Determine ranking process and who is responsible for each ranking
- Determine process to convert ranking results into risk reduction strategies and budget decisions

#### (2) Comparative Risk Analysis

- Finalize list of potential risks for ranking
- Gather data
- Assess residual risks and future risks for hazards to human health, ecosystems, and quality of life
- Address transboundary effects of risks
- Prioritize risks by ranking them qualitatively or quantitatively depending on data availability/needs
- Document risk analysis method
- Identify areas of uncertainty requiring more research/data
- Identify environmental indicators that will help monitor risks in the future

#### (3) Risk Management

- Assess and rank risks for management by:
  - identifying the most serious/adverse environmental and/or health risks;
  - identifying community values concerning raks; and
  - ranking risks in order of seriousness and community values.
- Select risk management factors
- Determine risk reduction goals
- Revise priorities and reduce risks by:
  - deciding which risks to address;
  - developing action plans to reduce risks; and
  - developing ongoing monitoring programs to ensure that risks are reduced effectively.
- Propose action plan activities to reduce or prevent risk, a schedule, and measures of progress
- Develop actions to overcome barners
- Document action plans
- Establish process for repeating risk ranking or updating results

- Assess dose-response relationship:
  - cancer potency factors for carcinogens
  - reference dose and other maximum safe levels for non-carcinogens
- Assess magnitude, duration, and frequency of exposure:
  - -- identify significant exposure pathways and routes
  - -- identify sources, location, timing, and quantity of pollutants released
  - -- describe fate and transport of pollutants
  - -- estimate concentration of pollutants
  - -- define human exposures (identify exposed populations, calculate intakes)
  - -- describe and document uncertainties in the data
- Characterize potential risks (cancer and non-cancer):
  - -- individuals
  - populations
- Rank potential risks (cancer and non-cancer):
  - quantitative factors (e.g., cancer incidence, non-cancer hazard indices)
  - -- qualitative factors (severity of effects, quality of data)
  - risk ranking exists along spectrum ranging from purely judgmental to rigorously quantitative
- Combine cancer and non-cancer risks:
  - -- very general guidance for this step
  - -- non-quantitative or semi-quantitative, depending on judgment and data availability
  - group potential cancer and non-cancer health effects together in matrix with population data to determine three separate categories of relative risk, e.g., "catastrophic," "serious," or "adverse"
  - if desired, an overall health rating can be obtained by aggregating all the health-effect category rankings, assuming equal weight for all rankings
- (2) Comparing and assessing ecological risks
- Determine and define environmental problem:
  - isolate/partition problem by geographical area or ecosystem type
  - select criteria to evaluate the problem (e.g., area of impact, severity or reversibility of impact, uncertainty, or "value" of ecosystem)

## Analyze the problem:

- identify stressors (e.g., chemicals) and establish causal link between problem and ecological effects
- estimate exposures (e.g., estimate concentrations)
- -- characterize ecological effects
- Characterize risks (using narrative and/or numeric descriptors):
  - -- describe each problem area using common evaluative criteria (e.g., severity of impact is "low")
  - -- summarize overall risk across stressors or problem areas to specific geographical area or ecosystem type
  - -- aggregate risks across all geographical areas or ecosystem types in the study

## • Compare and rank risks:

-- compare ecological risks posed by different problem areas and rank them into broad categories of relative risk (professional judgement and consensus-building are integral to this process)

## (3) Quality-of-life assessments

- Identify impacts of pollution on society and determine community values and social concerns:
  - -- spiritual, cultural, aesthetic, religious, ethnic values, or concern for future generations
  - -- environmental justice for diverse populations and lifestyles
  - -- economic stability for natural-resource-intensive industries

Ensures that assessment process has broad public support and accurately represents public concerns

• Define evaluative criteria based on broadly shared public values for evaluating effects of pollution on quality of life (e.g., no. of people affected, reduced recreational opportunities).

# Collect and analyze data:

- surveys, questionnaires, census data, public meetings
- analytic methods (e.g., damage to materials, commercial harvest losses, health care costs, recreational losses, property-value losses, resource restoration costs, aesthetic damage)

- Characterize impacts for all problem areas:
  - long-term damage (e.g., ozone depletion in the future will increase health care costs and damage PVC plastics)
  - loss of natural resources (e.g., oil, minerals, wood)
  - loss of natural ecosystem services (e.g., protection from UV radiation by the ozone layer, purification of water by wetlands)
- Present findings and rank impacts to quality of life using qualitative and/or quantitative data:
  - -- establish integrated ranking of impacts
  - document process, which may require controversial analytic methods
  - -- combine qualitative descriptions of impacts with dollar damage estimates
- Analyze future environmental conditions (i.e., incorporate longer-term viewpoint into the assessment of environmental problems).

#### System evaluation

## Scope of coverage:

- Selects problems from a core list of environmental problems (e.g., industrial wastewater discharges to oceans, lakes, and rivers; physical degradation of terrestrial ecosystems/habitats).
- Examines risks from current pollution whether short-lived or present for centuries and effects of today's pollution on the future. Does not try to evaluate risks from tomorrow's pollution due to the speculative nature of predicting trends.

## Media addressed:

• Media are considered to be integrated, not discrete. System evaluates interrelated causes and effects of pollution to air, water, and land.

# Types of targeting criteria used:

- Uses a broad range of targeting criteria, key criteria are summarized in Exhibit 2.
- Human health and ecological risk; adverse effects on quality of life.
- Inadequacies of environmental regulations: risks created by uncoordinated government actions and risks created by programs that fail to address complex relationships among environmental problems.
- Economic viability, technological feasibility, and social equity.

## Data requirements:

- Qualitative and quantitative data on risks to human health and ecosystems developed or gathered by the Technical Work Groups using the system, depending on availability and quality of data.
- Public ethics, values, and concerns (e.g., health of children, protection of ecosystems, wellbeing of future generations) as expressed by public forums and surveys.
- Use of professional judgment varies considerably during the individual risk rankings, depending on the quality and availability of necessary data.

# Applicability to waste minimization targeting:

#### Pro

- Broad framework allows flexibility in implementing the approach.
- Integrates judgment and technical expertise with values and concerns expressed by the community to determine which risks are of priority. Eliminates public confusion caused by information that is obsolete, incomplete, or biased to serve a particular viewpoint.

#### Con

- System is labor intensive and politically charged. Substantial investment of money and time necessitates careful planning for the process as a whole, not in segments.
- Problems are considered by the system to have interrelated causes and interrelated effects.
   May be difficult to adapt the system to discrete problems like RCRA wastestreams at specific facilities.
- Because of the significant latitude in structuring and scoring the criteria, and in relying on expert judgement, may get different results in different applications.

Exhibit 2: Targeting Criteria Used in the Regional Comparative Risk Ranking System

TARGETING CRITERIA	·
Direct Risks	
Waste volume	
Waste/constituent toxicity Human toxicity Ecological toxicity	1,
Constituent concentration	1
Waste type	
Number of generators	
Waste management practices	1
Releases to environmental media	1
Potential for constituent transport	1
Potential for cross-media transfer	1
Potential for human exposure: environmental settings occupational settings	
Potential for ecological exposure	
Indirect or Acute Risks	
Ignitability, corrosivity, reactivity	
Ozone depletion potential	1
Global warming	1
Other Targeting Criteria	
Hazardous waste management capacity	
Technical/administrative feasibility	1
Permitting/enforcement factors	
Cost savings	1
Other factors  Quality of life	1

# OPPTS Existing Chemicals Screening Program<sup>5</sup>

## Purpose of the Methodology/System

- "To screen, establish testing requirements for, assess, and develop strategies for managing risks posed by chemicals currently in production or use."
- To make decisions on regulatory and nonregulatory actions to reduce or eliminate the possibility of harm to human health or the environment.
- Program recently amended to increase effectiveness of its risk management actions, to increase public involvement and public understanding of the risks of chemicals, to incorporate the concept of pollution prevention, and to integrate the program with agency-wide risk reduction priorities.

#### System Description

• Program composed of three stages: Risk Identification, Risk Evaluation, and Risk Management.

#### Risk Identification

- Program receives and reviews substantial amount of risk-related information on chemicals submitted by chemical manufacturers and other sources in response to TSCA Section 8; this information includes:
  - -- production, use, and exposure data from manufacturers required under Section 8(a);
  - -- section 8(c) records of "significant adverse reactions;"

"EPA's Existing Chemical Program: An Overview" [no date or source]

U.S. EPA. 1991. RIB How-to Guide: RMI Economic Reports. Regulatory Impacts
Branch, Economics and Technology Division, Office of Toxic Substances, U.S. Environmental
Protection Agency, Washington, D.C.

U.S. General Accounting Office. 1984. Report by the Comptroller General of the United States: EPA's Efforts to Identify and Control Harmful Chemicals in Use. Gaithersburg, MD. GAO/RCED-84-100.

U.S. EPA. 1986. Toxic Substances Control Act (TSCA): Report to Congress for Fiscal Year 1985. U.S. Environmental Protection Agency, Washington, D.C.

Contact: John Leitzke (202/260-3507)

<sup>&</sup>lt;sup>5</sup>Documentation reviewed:

- section 8(d) health and safety studies; and
- section 8(e) substantial risk notifications.
- Other information reviewed by the program includes:
  - ... Chemical Hazard Information Profiles;
  - -- Substitute Hazard Profiles:
  - -- National Toxicology Program Studies; and
  - monitoring studies, which are reviewed to obtain data on levels of human and environmental exposure to substances of concern (for example, the National Adipose Tissue Survey is reviewed to detect TSCA-related chemicals in human tissue; occupational exposure monitoring studies are reviewed to assess human exposure)

#### Risk Evaluation

- EPA determines nature and magnitude of risks by analyzing health and environmental effects data gathered during risk identification activities
- EPA performs chemical-specific risk assessment based on both toxicity and exposure (e.g., for formaldehyde, EPA evaluated the two largest populations exposed to this chemical, i.e., permanent press apparrel manufactures and mobile home owners) and decides on need for risk management activities

# Risk Management

Two phases: Risk Management One (RM1) and Risk Management Two (RM2)

#### RM1

- Based on qualitative risk evaluations (i.e., initial screening that relies on readily available
  data on potential toxicity and potential exposure) developed for each chemical candidate,
  the RM1 committee selects one of four options:
  - placing chemical on the Master Testing List (i.e., a priority testing list) because of significant data gaps relating to the chemical's hazard or exposure potential;
  - -- placing chemical on the Risk Reduction List if it is believed or known to pose significant risks;
  - placing chemical on the Regional Activities Track if concern for chemical is limited to specific geographic regions; or
  - -- dropping chemical from the list of candidates for risk management.
- From chemicals placed on Risk Reduction List, OPPTS selects a subset as potential candidates for action under RM2 phase, using four criteria:

- -- TSCA jurisdiction:
- potential or known toxicity;
- potential or known exposure to the chemical; and
- -- potential for pollution prevention.

(Only chemicals from the Risk Reduction List enter the RM2 stage.)

- The documentation reviewed did not describe what happens to chemicals that are not placed on the Risk Reduction List.
- Important point: From the documentation reviewed, it appears that there is no standard system or "cook-book" type approach by which each chemical is evaluated in the risk management process. Rather, the priority of action for each chemical in the Existing Chemicals Program (i.e., a chemical's relative "riskiness") is determined on a case-by-case basis through a review of risk assessments and all relevant information gathered, as discussed above, and through coordination among various divisions in OPPTS. This review for each chemical in the RM1 phase takes the form of a 12-week review cycle, approximating the steps outlined in Exhibit 1.

# Exhibit 1. The RM1 Project Cycle

- 1. The Existing Chemical Assessment Division (ECAD) coordinates the initial screening process to select chemical candidates for the RM1 phase. ECAD prepares a dossier on the chemical, and, along with three other OPPTS divisions involved in the project cycle, has a schedule in which to complete its assigned report that will make up the dossier.
- 2. The Economics and Technology Division/Industrial Chemistry Branch prepares a chemistry report for the chemical under review during Week 1.
- 3. The Economics and Technology Division/Regulatory Impacts Branch prepares the RM1 economics report during Week 3. The report is based on data elements that include volume information, market trends, use data, producer data, importer data, and substitutes data.
- 4. The Economics and Technology Division/Chemical Engineering Branch prepares the engineering report which evaluates worker exposures and releases into the environment for the chemical during Week 5.
- 5. The Health and Environmental Review Division (HERD), during Week 5, prepares the hazard report which reviews the health and environmental hazards posed by the chemical under review.
- 6. The Exposure Evaluation Division/Exposure Assessment Branch, during Week 7, prepares the exposure report which evaluates the environmental concentrations of the chemical as well as consumer and general public exposures.
- 7. The Existing Chemical Assessment Division prepares a dossier based on the reports listed above during Week 9.

#### RM2

- Existing information is investigated thoroughly so that OPPTS may better its understanding of the chemical's hazard and risks and, more specifically, determine which portions of a chemical's lifecycle (manufacture, processing, distribution, use, or disposal) pose risks.
- Results of the investigation are used to select a risk management strategy that may consist of multiple components including:
  - -- public awareness campaign;

- call to industry for voluntary action;
- enforcement of existing regulations;
- -- development of new regulations (e.g., bans, labeling requirements); and
- -- removal of the chemical from further consideration.
- Potential risk reduction actions are developed, emphasizing source reduction, responsible recycling, improved treatment technologies, and improved disposal technologies.

#### System Evaluation

## Scope of coverage:

- Chemical-specific; does not consider wastes or wastestreams specifically.
- Population of potential chemical candidates evaluated in the initial screening of the RM1
  phase consists of the approximately 14,000 chemicals found on the TSCA Inventory whose
  annual production quantities exceed 10,000 pounds.
- Considers risk from chemical's entire lifecycle (manufacture through disposal)

## Media addressed:

 Because program is a case-by-case evaluation of risks posed throughout a chemical's lifecycle, it appears that multiple media and exposure pathways would be considered (although they may differ from chemical to chemical).

# Types of targeting criteria used:

- Statutory: Language in TSCA gives EPA discretion to determine "unreasonable risk" posed by a chemical to human health or the environment; such a determination is the trigger mechanism for control action under TSCA. In assessing unreasonable risk, EPA must consider the following criteria in initial RM1 screening activities:
  - -- human health and environmental hazard;
  - -- degree of human and environmental exposure;
  - the benefits provided by the chemical's uses;
  - the availability of substitutes for such uses; and
  - the economic consequences of regulating the chemical, considering impacts on national economy, small business, technological innovation, the environment, and public health.
- RM2: The four criteria used by OPPTS in moving chemicals from the Risk Reduction List to the RM2 phase as candidates for risk management activities are:
  - TSCA jurisdiction
  - potential or known toxicity of the chemical

- potential or known exposure to the chemical
- potential for pollution prevention
- Targeting criteria used in the Existing Chemicals Program are summarized in Exhibit 2.

#### Data Requirements:

- Initial screening of chemicals prior to the RM1 stage utilizes TSCA section 4 test rule data:
  - EPA has authority under Section 4 of TSCA to require by rule that chemical manufacturers and processors test for various health and environmental effects.
  - -- test rules imposed on industry by EPA:
    - are chemical-specific;
    - require testing for specific human health and environmental effects;
    - specify test standards;
    - impose submission deadlines for test data; and
    - assign responsibility for testing.
- Based on reviewed documentation, it appears that the reporting process in the RM1 project cycle is data-intensive and requires the various OPPTS divisions to supplement existing information in order to fulfill certain data requirements.
  - example: Regulatory Impacts Branch (RIB), in preparing the economics report, has to base its research on readily available secondary sources and must conduct a library search for data on the volume, market trends, uses, producers and importers, and substitutes for the RM1 chemical; examples of sources include:
    - SRI. Chemical Economics Handbook
    - SRI, Directory of Chemical Producers
    - EPA Regulatory Analyses
    - on-line search of chemical industry business and chemistry data bases available through DIALOG and STN International

# Applicability to Waste Minimization Targeting:

#### Pros

- Several stages of the program incorporate the concept of pollution prevention (although from the documentation available, it's not clear what 'subcriteria' are used to determine a chemical's potential for pollution prevention).
- The chemical targeting system is intertwined with a system of risk management strategy development.

#### Cons

- The program is designed to address individual chemicals and not wastes, wastestreams, or industrial establishments.
- The program has no standard system of risk evaluation; rather, chemicals are evaluated on a case-by-case basis by various branches/divisions within OPPTS.
- The preparation of the chemical dossiers involves coordination among many branches and divisions of OPPTS and is time- and resource-intensive. This approach accounts for many relevant factors and develops consensus on the 'riskiest' chemicals.

Exhibit 2. Targeting Criteria Used in the OPPTS Existing Chemicals Screening Program

TARGETING CRITERIA						
Direct Risks						
Waste volume						
Waste/constituent toxicity  Human toxicity  Ecological toxicity	1					
Constituent concentration or mass	1					
Waste type (e.g., solvents)	•					
Number of generators	1					
Waste management practices						
Releases to environmental media	1					
Potential for constituent transport						
Potential for cross-media transfer						
Potential for human exposure: residential settings occupational settings	4					
Potential for ecological exposure	1					
Indirect or Acute Risks						
Ignitability, corrosivity, reactivity						
Ozone depletion potential						
Global warming						
Other Targeting Criteria						
Hazardous waste management capacity						
Technical/administrative feasibility						
Permitting/enforcement factors						
Cost savings						
Other factors  Potential for pollution prevention  Availability of substitutes  Economic impacts of regulating chemical  Benefits of chemical uses	***					

## Industrial Pollution Prevention Opportunities for the 1990s6

#### Purpose of the Methodology/System

- System was a one-time research effort to identify a short list of industries, or industry segments, that present significant opportunities for waste reduction and pose environmental problems or risks in terms of the wastes generated.
- More generally, the system was expected to "provide a data base that could be used as guidance by the EPA for the development of a research strategy for pollution prevention."

#### System Description

- System is based on subjective, expert evaluation of 12 criteria (see Exhibit 1). Of the 12, 11 are non-risk-based criteria (listed under "Other factors" in Exhibit 1).
- First, a list of 175 industries was selected from the 1987 SIC publication based on the selection criteria.
- Second, the 175 SIC-based industries/industrial sectors were shortlisted to a group of 20 by experts from USEPA, academia, state pollution prevention programs, and contractor personnel, again using the 12 selection criteria. The 20 SICs were ranked in priority order by the experts.
- Finally, the experts "subjectively normalized" the SIC-based industries (aggregating the multi-segmented SICs) and developed a list of 17 industry segments (see Exhibit 2) for further research into pollution prevention opportunities. This list is supposed to best represent the problems and opportunities for pollution prevention.

#### System Evaluation

## Scope of coverage:

• Covers industries/industrial sectors covered by four-digit SIC codes; system does not use information at the wastestream or constituent level.

## Media addressed:

System does not explicitly consider releases or threat of releases to any media; experts may or may not have considered such releases when selecting the priority industries.

Contact: Ivars Licis (513/569-7718)

<sup>6</sup> Documentation reviewed:

U.S. EPA. Aug. 1991. Industrial Pollution Prevention Opportunities for the 1990s. Prepared by the Office of Research and Development, U.S. Environmental Protection Agency, Washington, D.C., EPA/600/8-91/052.

Exhibit 1: Targeting Criteria Used in the Industrial Pollution Prevention Opportunities for the 1990s System (cont.)

TARGETING CRITERIA	
Direct Risks	
Waste volume (pounds)	1
Waste/constituent toxicity	1
Constituent concentration	
Waste type	1
Number of generators	
Waste management practices	
Releases to environmental media	
Potential for constituent transport	
Potential for cross-media transfer	
Potential for human exposure: environmental settings occupational settings	
Potential for ecological exposure	
Indirect or Acute Risks	
Ignitability, corrosivity, reactivity	
Ozone depletion potential	
Global warming	
Other Targeting Criteria	
Hazardous waste management capacity	
Technical/administrative feasibility	
Permitting/enforcement factors	
Cost savings	

Exhibit 1: Targeting Criteria Used in the Industrial Poliution Prevention Opportunities for the 1990s System (cont.)

Other factors	·
Importance of industry to society or nation	1
Large frequency of small to mid-size firms that would benefit from govt. participation	1
Significant benefits would be derived from WM	1
WM would not adversely impact product quality or marketability	1
WM would offer cost benefits in the long run	1
WM in this industry would be readily transferable to others	
Industry has exhibited interest in WM	
WM appears to be technologically achievable	1
Industry would benefit from govt. involvement	1
Industry would be receptive to WM studies	1
Industry will not be viable in the long run without massive changes	•

## Types of targeting criteria used:

- Targeting criteria used in this system are varied; some relate, at least indirectly, to risk, and others to measures such as the industry's potential for success in pollution prevention.
- Targeting criteria were chosen for evaluating or comparing the relative importance, as perceived by an expert, of certain industry characteristics/attributes that would relate to the industry's feasibility for pollution prevention research. Examples of these attributes include industry size, waste production in terms of toxicity and/or volume, and receptivity of the industry to innovation.
- System relied on professional judgement for factors such as the degree to which each criterion had to be satisfied and the number of criteria that had to be satisfied simultaneously by the industry being reviewed.

## Data requirements:

• Documentation does not explain what data were used by the experts in selecting the 17 priority industry segments; only "best informed judgement" is mentioned.

# Applicability to Waste Minimization Targeting:

#### Pros

• 17 industry segments that have expert consensus on their feasibility for pollution prevention can be used as starting point for identifying specific wastestreams for WM.

Cons

- Based on available documentation, methodology used to prioritize industries is not well defined and does not appear to reproducible.
- System cannot be used again without reconvening an expert panel.
- Neither specific targeting criteria nor data needs are well defined.

Exhibit 2: Priority Industry Segments Identified for Pollution Prevention Research 1

Textile dyes and dyeing
Wood preserving
Pulp and paper
Printing
Chemical manufacture
Plastics
Pharmaceutical
Paint industry
Ink manufacture
Petroleum industry
Steel industry
Non-ferrous metals
Metal finishing
Electronics/semiconductors
Automobile manufacture/assembly
Laundries/dry cleaning
Automobile refinishing/repair

<sup>&</sup>lt;sup>1</sup> Because they are "normalized" from a larger set of industry segments, these 17 do not directly correspond to any SIC.

# National Corrective Action Prioritization System (NCAPS)7

# Purpose of the Methodology/System

- Used to prioritize treatment, storage, and disposal facilities (TSDFs) regulated under RCRA for possible corrective actions.
- Meant to be an internal EPA management tool and is not subject to external review, comment, or approval
- Incorporates many of the same factors and equations as the Hazard Ranking System (HRS), but is much simpler and less data-intensive. The equations for pathway scores are conceptually similar to those in the HRS in that they involve multiplication of release, waste characterization, and targets related factors, and division by a scaling factor. However, the factors are scored in a much simpler fashion.

#### System Description

- NCAPS facility migration score determined by evaluating four routes of potential contamination:
  - -- ground water
  - -- surface water
  - -- air
  - -- on-site exposure
- Each route of potential contamination scored based on the scores assigned to specific factors (illustrated in Exhibit 1) that fall into one of the five following categories:
  - -- releases
  - -- route characteristics
  - -- containment
  - -- waste characteristics
  - targets
- Factor scores are combined using route-specific and scenario-specific formulae (e.g., formulae dependent on the presence of an observed release or a possible release) and the resulting values are normalized to generate route scores between 0 and 100.

U.S. EPA. 1991. Environmental Protection Agency Technical Enforcement Support at Hazardous Waste Sites: National RCRA Corrective Action Prioritization System Guidelines Revised. Prepared by ICAIR Life Systems, Inc. Revised by PRC Environmental Management, Inc.

RCRA Prioritization System Scoring Summary [no other citation information].

Contact: Dave Fagan (703/308-8620)

<sup>&</sup>lt;sup>7</sup>Documentation reviewed:

- it appears that the scoring process for each route is strongly dependent on the presence of an observed release.
- Facility migration score is calculated by combining the four route scores in a root-mean-square equation.
- Chemical- or waste code-specific data requirements used in the scoring process include:
  - -- toxicity, based on the Sax rating (i.e., a toxicity rating scheme, that emphasizes acute toxicity, used in the original HRS)
  - -- persistence
  - -- waste quantity (based on known or estimated amounts)
- Concentrations of contaminants are <u>not</u> considered. However, waste <u>quantity</u> is considered (e.g., for scoring observed releases, the amount of waste actually released is considered).
- Factors in the releases, route characteristics, containment, and targets categories are generally exposure-related.

#### System Evaluation

# Scope of coverage:

- NCAPS scores and ranks TSDFs covered under RCRA.
- RCRA constituents and wastes.

#### Media addressed:

- Ground water
- Surface water
- Air
- Direct on-site exposure (i.e., direct physical contact):
  - nearby residential population with potential access to a site
  - -- sensitive environmental populations

#### Types of targeting criteria used:

- In general, NCAPS scoring factors are risk-based, relating to exposure (e.g., observed and possible releases, route characteristics factors, targets factors, persistence, and containment) and toxicity Key targeting criteria are summarized in Exhibit 2.
- Some NCAPS targeting factors are based on permit information. For example:

- scoring for releases in the surface water route is partially dependent on whether the discharge (outfall) was permitted and whether permit violations have occurred at the facility.
- in such cases, scoring the <u>observed</u> release factor applies only to facilities with unpermitted discharges while scoring the <u>possible</u> release factor applies to facilities with permits.
- Facility history may also affect NCAPS scoring. For example:
  - for facilities that have removed their hazardous wastes but have past containment practices ranked less than 'very good', scoring for releases uses the past containment rating
  - the air route is scored for the possible release factor if residents have complained of odors or if a facility investigator has noted odors

#### Data requirements:

- Not especially data-intensive (relative to HRS)
- In the absence of site-specific information for scoring some factors, the NCAPS guidance supplies default assumptions that can satisfy the data requirements of the scoring equations.
- Types of data sources that can be used in the NCAPS scoring process:
  - -- documentation of releases of hazardous waste or hazardous constituents:
    - RCRA Facility Assessment (RFA)
    - Preliminary Review/Visual Site Inspection (PR/VSI)
    - CERCLA Preliminary Assessment (PA) report
    - Preliminary Assessment/Visual Site Inspection (PA/VSI) report
  - -- waste quantity information:
    - RCRA Part A permit application
    - tank capacity
    - permitted drum storage capacity
    - other site report information
    - targets:
      - quantification of population size is not required
      - wetlands, streams, rivers, and residential areas as well as their distances from the facility can be identified in a United States Geological Survey map

# Applicability to waste minimization targeting:

#### Pros

Considers RCRA wastes and RCRA constituents.

Factors used for scoring conform to risk assessment requirements.

- Unlike other systems such as the HRS, NCAPS does not carry the weight of external review, comment, or approval
- May not be sufficiently quantitative; for example:
  - -- population size potentially affected by releases is not quantified
  - -- waste characteristics include only waste quantity, toxicity, and persistence; several other characteristics are not considered, such as bioaccumulation potential and contaminant concentrations
  - incorporates multiple assumptions that may be used in the absence of specific information (for example, when evaluating facilities in communities adjacent to any Great Lake, the communities are assumed to obtain their residential water supply from the lake unless there are available data to disprove this assumption)
- In scoring an entire facility, NCAPS does consider releases from individual active units, which could include combustion units.

Exhibit 1. Factor Categories and Factors Used in NCAPs Scoring

Factor Category	) actors
Releases to the Environment	Ground water route observed releases Surface water route observed releases presence of permitted outfalls presence of permit violations Air route presence of observed, unpermitted, ongoing releases existence of facility operating air permit permit violations or odor complaints by residents can contaminants migrate into air? containment
Route Characteristics (evaluated when an observed release has not been scored for a particular route)	Ground water route depth to aquifer net precipitation physical state Surface water route facility location 24-hour rainfall distance to surface water physical state On-site exposure route site accessibility
Containment	Ground water route evaluates containment properties of: surface impoundments containers/tanks piles landfills Surface water and on-site exposure routes evaluate containment properties of: surface impoundments containers/tanks piles landfills Air route evaluates containment properties of: closed and open containers and tanks
Waste Characteristics	Toxicity Persistence Waste quantity

# Exhibit 1 (continued). Factor Categorias and Factors Used in NCAPs Scoring

Factor Category	Factors
Targets	Ground water route ground water use distance to intake Surface water route surface water use distance to intake or contact point distance to sensitive environment Air route population (residences, industries, agricultural lands, other) distance to sensitive environments On-site exposure route distance to residential areas on-site sensitive environments

Exhibit 2. Targeting Criteria Used in the National Corrective Action Prioritization System

TARGETING CRITERIA		
Direct Risks		
Waste volume	1	
Waste/constituent toxicity Human toxicity Ecological toxicity	1	
Constituent concentration or mass	✓ (mass only)	
Waste type (e.g., solvents)	1	
Number of generators	·	
Waste management practices		
Releases to environmental media	1	
Potential for constituent transport		
Potential for cross-media transfer	1	
Potential for human exposure: residential settings occupational settings	1	
Potential for ecological exposure	1	
Indirect or Acute Risks		
Ignitability, corrosivity, reactivity		
Ozone depletion potential		
Global warming		
Other Targeting Criteria		
Hazardous waste management capacity	1	
Technical/administrative feasibility		
Permitting/enforcement factors	1	
Cost savings		
Other factors Odor complaints by residents Exposure to agricultural lands	1	

# Minnesota Office of Waste Management Nonhazardous Industrial Waste Targeting and Pollution Prevention Project<sup>2</sup>

#### Purpose of the Methodology/System

- Identify high-risk nonhazardous industrial wastes to target them for pollution prevention demonstration projects.
- Results from demonstration projects will be used to develop fact sheets and case study summaries to assist industry in achieving additional pollution prevention.

#### System Description

- Two hazard ranking (HR) values are developed for nonhazardous wastes managed in landfills: a Toxicity Characteristic Leaching Procedure (TCLP) value and a Recommended Allowable Limit (RAL) value. (RALs are established by the Minnesota Department of Health as standards for safe levels of chemicals in drinking water from private wells.)
  - 1. TCLP HR = annual waste volume \* Σ (measured TCLP level/TCLP regulatory level for constituent i)
  - 2. RAL HR = annual waste volume \* Σ (measured TCLP level/RAL for constituent i)
- Targeting is based on the TCLP HR and the judged potential for successful pollution prevention demonstration projects. (Documentation for this system does not explain how the RAL HR is used.)

#### System Evaluation

#### Scope of coverage:

- Industrial nonhazardous wastes managed in landfills, excluding one-time wastes (e.g., from spill cleanups).
- TCLP constituents.

Nonhazardous Industrial Waste Targeting and Pollution Prevention Demonstration Project, Interim Report, undated, prepared by the Minnesota Office of Waste Management under USEPA Solid Waste Management Assistance Grant Project X819717-01-0.

<sup>&</sup>lt;sup>8</sup> Documentation reviewed:

# Media/pathways addressed:

Ground water migration pathway.

# Types of targeting criteria used (see Exhibit 1):

- Risk based: Waste volume plus all of the fate, transport, exposure, and toxicity factors considered in the development of TC regulatory levels.
- Non-risk based: Potential for successful pollution prevention; extent to which pollution prevention demonstration projects have already been conducted.

#### Data requirements:

- Waste volume: volume managed in landfills.
- Waste/constituent toxicity: TC regulatory levels; Minnesota RALs in drinking water wells.
- Constituent concentration (in leachate): TCLP results.
- Data are derived from (1) municipal/industrial waste co-disposal applications filed by waste generators; (2) annual reports filed by co-disposal facilities; and (3) data collected by a commercial nonhazardous waste landfill.

# Applicability to waste minimization targeting:

#### Pros

- Calculations are simple and straightforward.
- Only two data elements are needed.

- Applicable only to landfilled waste (methodology can probably be adapted to other land management practices, but not to incineration).
- TCLP data may not be available for certain wastes/constituents, or, if available, may not be easily accessible if methodology is to be implemented by Headquarters.
- Method is designed for nonhazardous wastes but can be used for hazardous wastes if data are available.
- Does not account for additional risk via other exposure pathways, risk due to constituents other than TCLP constituents, or risk to ecological receptors.

Exhibit 1. Targeting Criteria Used in the Minnesota Office of Waste Management
Nonhazardous Industrial Waste Targeting and Pollution Prevention Project.

TARGETING CRITERIA		
Direct Risks		
Waste volume	1	
Waste/constituent toxicity Human toxicity Ecological toxicity	1	
Constituent concentration or mass	1	
Waste type (e.g., solvents)		
Number of generators		
Waste management practices		
Releases to environmental media	<i>,</i> a	
Potential for constituent transport	<i>A</i>	
Potential for cross-media transfer		
Potential for human exposure: residential settings occupational settings	,,	
Potential for ecological exposure		
Indirect or Acute Risks		
Ignitability, corrosivity, reactivity		
Ozone depletion potential		
Global warming		
Other Targeting Criteria		
Hazardous waste management capacity		
Technical/administrative feasibility	1	
Permitting/enforcement factors		
Cost savings		
Other factors		

These are considered indirectly in deriving the measured and regulatory TC levels.

# Numerical Hazard Ranking Scheme for Waste Scheduling

# Purpose of the Methodology/System

- Numerical scheme developed for ranking, according to hazard, the hazardous wastes listed in 40 CFR Part 261.
- At the time of development, intended to be used by EPA for scheduling the review of listed wastes with regard to land disposal restrictions.

# System Description

- Relatively ranks wastestreams based on a "Hazard Index" (HI) score; the HI score (1 to 100) is calculated as the product of the wastestream's toxicity and exposure scores.
- Wastestreams with multiple constituents are ranked based on the toxicity and exposure scores of the single highest scoring constituent.

# Toxicity Score:

- Scored from 1 to 9 based on chronic toxicity (using acceptable daily intakes or ADIs) or on carcinogenicity (using unit carcinogenic risk or UCR). The chronic toxicity score is adjusted upward by 1 if the constituent scores "high" for acute toxicity (based on LD<sub>50</sub>, LC<sub>50</sub>, or LD<sub>10</sub>).
- All constituents are scored for toxicity; system provides a data hierarchy to score all constituents, irrespective of how well they are characterized in terms of toxicity.

# Exposure Score:

Scored from 1 to 10 based on a total exposure dose, where

- Exposure dose calculated for a receptor under a <u>fixed</u> set of environmental conditions, i.e., releases from a landfill reaching the receptor located 100 meters from the source.
- Exposure dose calculated in three steps: (1) calculation of release rates from landfill to air, ground water, and surface water; (2) fate and transport modeling of steady-state concentrations at the exposure point; and (3) calculation of media-

<sup>&</sup>lt;sup>9</sup> Documentation reviewed:

Draft Interim Report on the Development of Numerical Hazard Ranking Scheme. April 19, 1984. Developed for the Office of Solid Waste by Environ Corp.

specific human intakes. (The surface water intake is estimated based on bioaccumulation in fish that is ingested.)

#### System Evaluation

#### Scope of coverage:

- System was applied to all <u>listed</u> hazardous wastestreams in 40 CFR Part 261, i.e., F, K, U, and P wastes.
- A composition profile was developed for each wastestream based on only those constituents in Appendix VII of 40 CFR Part 261.
- Documentation notes that the system was applied to approximately 450 wastestreams.

#### Media addressed:

 System is designed to provide rankings based on releases to air, surface water, and ground water.

#### Types of targeting criteria used:

• Targeting criteria include initial concentration of wastestream constituents and their toxicity. Initial concentrations dictate the release into various environmental media, and hence, the ultimate exposure dose. (See also Exhibit 1.)

#### Data requirements:

- System documentation provides a fair amount of detail on data sources, and methods for estimating some of the required parameter values.
- For waste composition information, system relied heavily on (1) data compiled for EPA's Risk-Cost Analysis Model and other "best available data" (for F and K wastes); and (2) assumptions about "typical upperbound concentrations at which commercial chemical products may become wastes" (for U and P wastes).
- For toxicity information, system relied first on available EPA ADIs and UCRs, and then on open literature, using human and animal toxicity data to derive ADIs and UCRs.

# Applicability to waste minimization targeting:

#### Pros

System directly applicable to RCRA wastestreams.

- Ranking is based on a fixed set of exposure conditions (i.e., assumptions regarding releases from a landfill and the location of the receptor).
- Uses certain conservative or "worst-case" simplifying assumptions for the fate and transport modeling (e.g., constant wind direction towards receptor, no decay of constituents in surface water).
- May not be as applicable for wastes destined for incineration because rankings appear to be driven by the ground-water pathway.
- Limited in targeting criteria, e.g., does not consider ecological toxicity, waste volume.

Exhibit 1. Targeting Criteria Used in the Waste Scheduling Scheme.

TARGETING CRITERIA		
Direct Risks		
Waste volume		
Waste, constituent toxicity  Human toxicity  Ecological toxicity		
Constituent concentration or mass	1	
Waste type (e.g., solvents)	•	
Number of generators		
Waste management practices		
Releases to environmental media (e.g., simulated releases from waste management units)	\$	
Potential for constituent transport (based on physico-chemical properties)	\$	
Potential for cross-media transfer		
Potential for human exposure: residential settings occupational settings	\$	
Potential for ecological exposure		
Indirect or Acute Risks		
Ignitability, corrosivity, reactivity		
Ozone depletion potential		
Global warming		
Other Targeting Criteria		
Hazardous waste management capacity		
Technical/administrative feasibility		
Permitting/enforcement factors		
Cost savings		
Other factors		

These are considered indirectly in modeling the exposure dose.

# Risk-Based Enforcement Strategy (RBES)<sup>10</sup>

#### Purpose of the Methodology/System

• To rank sites for enforcement activities and assess the effectiveness of environmental laws in reducing risks at these sites.

#### System Description

- System developed through a cooperative effort between the Office of Research and Development (ORD) and the Office of Enforcement (OE).
- Ranks facilities based on a Chemical Ranking Factor (CRF), compliance history, and environmental vulnerability.
- CRF is essentially the summation of the toxicity-weighted annual chemical releases at a facility. (EPA selected this over a more complicated approach to calculate the CRF whi instead relied on exposure-adjusted releases.)
- To score the CRF, system f. ... selects facility or industrial sector to be ranked, and retrieves release information, on a facility-specific basis, from three EPA information systems:
  - Aerometric Information Retrieval System's Facility Subsystem (AFS-AIRS) a data base that contains data on major industrial, commercial, and municipal facilities required to report their emissions to EPA under the Clean Air Act. AFS-AIRS contains data primarily from SIC code 4911 (coal-burning and other power plants).
  - Permit Compliance System (PCS) an inventory for the National Pollutant Discharge Elimination (NPDES) program that includes permit and compliance information on regulated facilities (i.e., facilities discharging wastewater into navigable waterways). PCS contains data primarily from SIC code 4952 (sewage systems).
  - Toxic Release Inventory (TRI) contains data relating to manufacturing (SIC codes 2000-3999); facilities with at least 10 employees and that manufacture, process, or import at least 25,000 pounds of a TRI-listed chemical must file TRI reports. TRI primarily contains data from SIC codes 2000-3999 (manufacturing).

These three information systems contain data about three different types of facilities; however, there is some overlap. In using the system, total TRI water and total TRI air releases are compared to those from PCS and AFS-AIRS, respectively, and the larger value for each media is selected.

<sup>10</sup> Documentation reviewed:

Risk-Based Enforcement Strategy II, Draft Report Submitted to the Exposure Assessment Group, Office of Health and Environmental Assessment, U.S. EPA by Versar, Inc., September 30, 1993.

- Second, the release data is linked to a toxicity data base, and each chemical is assigned a chemical toxicity score equal to the inverse of its reportable quantity (RQ).
- Third, each chemical release is multiplied by the chemical-specific toxicity score; chemical scores for a given media are summed to yield a media score.
- Finally, all media scores for a facility are summed to yield a facility score.

#### System Evaluation

#### Scope of coverage:

- Covers all facilities for which release information exists in the AFS-AIRS, PCS, and TRI national data bases.
- Covers chemicals that have EPA-assigned RQs, or have sufficient toxicity information that an RQ could be derived for purposes of this system.

#### Media addressed:

System is designed to provide rankings based on releases to air, surface water, and land, both for on-site releases and off-site transfers

#### Types of targeting criteria used:

Targeting criteria include annual chemical releases at facilities and the toxicity of the released chemicals, both of which are incorporated into the CRF. (Although "compliance history" and "environmental vulnerability" are also mentioned in the documentation as facility-level criteria, no mention is made of how they are incorporated into the final facility ranking.)

# Data requirements:

- Uses release information available in the AFS-AIRS, PCS, and TRI national data bases.
- Uses chemical toxicity information stored in the toxicity data base developed specifically for this system.

# Applicability to waste minimization targeting:

#### Pros

• System may be applied to sources such as hazardous waste sites, waste disposal facilities, or manufacturing facilities; uses release information from numerous data bases that cover several different types of facilities.

- System appears relatively new, not well documented, and not sufficiently tested.
- Relies on only toxicity and annual releases and does not directly address exposure.

Exhibit 1. Targeting Criteria Used in the Risk-Based Enforcement Strategy.

TARGETING CRITERIA		
Direct Risks		
Waste volume		
Waste/constituent toxicity Human toxicity Ecological toxicity	1	
Constituent concentration or mass	1	
Waste type (e.g., solvents)		
Number of generators		
Waste management practices		
Releases to environmental media	1	
Potential for constituent transport		
Potential for cross-media transfer		
Potential for human exposure: residential settings occupational settings		
Potential for ecological exposure		
Indirect or Acute Risks		
Ignitability, corrosivity, reactivity		
Ozone depletion potential		
Global warming		
Other Targeting Criteria		
Hazardous waste management capacity		
Technical/administrative feasibility		
Permitting/enforcement factors		
Cost savings		
Other factors		

# Superfund Hazard Ranking System<sup>11</sup>

#### Purpose of the Methodology/System

- Used by EPA's Superfund program to assess the relative threat associated with actual or potential releases of hazardous substances.
- Primary screening tool for determining whether a site is to be included on the National Priorities List (NPL), which contains EPA's priorities for further investigation and possible remedial response action under Comprehensive Environmental Response, Compensation, and Liability Act (CERCLA).

#### System Description

- HRS site score determined by evaluating four pathways:
  - -- ground water migration
  - -- surface water migration
  - -- air migration
  - -- soil exposure
- Each pathway score based on a number of subcriteria, called 'factors', grouped into three primary criteria called 'factor categories':
  - likelihood of release (for the soil exposure pathway, likelihood of exposure);
  - -- waste characteristics: and
  - -- targets (e.g., the people or sensitive environments affected by the release).
- Factor category scores are determined from evaluation of factors. They are multiplied together and then normalized to 100 points to obtain pathway score.
- Factors with larger associated relative threats may carry greater weight and will more strongly affect final scores (for example, in the targets factor category, the HRS gives

Contact: Janet Grubbs (703/603-8833)

<sup>&</sup>lt;sup>11</sup>Documentation reviewed:

U.S. EPA. 1992. Hazard Ranking System Guidance Manual. Hazardous Site Evaluation Division, Office of Solid Waste and Emergency Response, U.S. Environmental Protection Agency, Washington, D.C., EPA 540-R-92-026.

U.S. EPA. 1990. The Revised Hazard Ranking System: Background Information. Hazardous Site Evaluation Division, Office of Solid Waste and Emergency Response, U.S. Environmental Protection Agency, Washington, D.C., Publication 9320.7-03FS.

greater weight to actual exposures, such as to people whose drinking water is contaminated and to actual contamination of the aquatic human food chain).

- Site score is obtained by combining the four pathway scores in a root-mean-square equation:
  - -- root-mean-square approach gives more weight to higher scoring pathways
  - -- scores range from 0 to 100
  - -- score of 28.50 or greater qualifies site for the NPL.
- Exhibit 1 is a summary of the main criteria and subcriteria (i.e., factor categories and factors) considered in the four HRS pathways, and is organized by factor category.
- Factors in the HRS utilizing chemical-specific data in the scoring process:
  - -- human toxicity factor (in the waste characteristics factor category), intended to represent the relative potential of a substance to cause adverse health effects, based on three measures of toxicity in a tired approach:
    - cancer, based on cancer potency factors and weight-of-evidence (ED<sub>10</sub> can also be used
    - noncancer effects of chronic exposure, based on verified Reference Doses (RfDs)
    - acute toxicity, based on LD50/LC50 data
  - hazardous waste quantity factor (in the waste characteristics factor category) based on hazardous constituent concentration data, mass of waste as deposited, volume, and/or surface area of the source.
  - -- several other chemical-specific waste characteristics factors include: persistence, mobility, bioaccumulation, ecological toxicity
- Almost all factors in the likelihood of release and targets factor categories are exposurerelated. For example:
  - -- observed release factor (in the likelihood of release factor category) based on a measured concentration of a hazardous substance in the environment to which a population is potentially exposed
  - targets factor category utilizes factors related to actual population exposures, which are depicted in Exhibit 1.

#### System Evaluation

Scope of Coverage:

- Constituents: HRS considers CERCLA hazardous substances as well as other pollutants or contaminants:
  - -- hazardous substances defined in CERCLA section 101(14), which references substances specifically listed under other Federal laws.
  - "pollutants or contaminants" broadly defined in CERCLA section 101(33) and could include any constituent reasonably anticipated to be harmful to human or ecological health; EPA determines on a case-by-case basis which substances fall within definition.
  - -- available chemical data bases used in scoring (Superfund Chemical Data Matrix) focus on approximately 300 commonly encountered hazardous substances
- Sources: HRS considers "any area where a hazardous substance has been deposited, stored, disposed, or placed, plus those soils that may have become contaminated from hazardous substance migration", such as:
  - -- above-ground and below-ground tanks
  - -- contaminated soil (excluding land treatment)
  - -- drums
  - -- landfarms/land treatment
  - -- surface impoundments

#### Media/pathways addressed:

- Ground water migration pathway
- Surface water migration pathway (overland/flood and ground water to surface water components)—divided into three threats: drinking water, human food chain, and environmental
- Soil exposure pathway—divided into two threats: resident population and the nearby population
- Air migration pathway

# Types of targeting criteria used:

• The HRS scores a site using criteria based on human, environmental, and resource risk in multiple pathways. Key criteria in the HRS are summarized in Exhibit 2.

#### Data Requirements

• All available site information, which may include information from Preliminary Assessment and Site Inspection reports, should be collected by scorer and should include data on:

- primary sources or hazardous substances at the site
- the hazardous substances themselves and their quantities
- whether there are observed releases
- -- major targets (e.g., populations, municipal wells, fisheries, sensitive environments) located near the site
- -- whether any targets are exposed to actual contamination.
- Scorer should assess whether available information is sufficient to document all the HRS factors relevant to the site's score; additional data collection may be necessary to gain a better understanding of those factors <u>critical</u> to the site's HRS score.
- HRS accommodates various levels of data quality, and often provides default options when complete data is not available.
- Scorer should develop selective scoring strategy because it is usually not feasible to gather data for and score every factor in every pathway, and should consider the following:
  - a primary goal of HRS scoring is to determine whether or not the site is eligible for the NPL, i.e., if site score is greater than 28.50
  - -- sites often pose significant threats in only one or two pathways
  - higher-scoring pathways exert a proportionately greater influence on the site score than do lower-scoring pathways.

# Applicability to Waste Minimization Targeting:

#### Pros

- HRS is comprehensive, e.g., addresses numerous criteria in four pathways, and thus may accurately assess threats associated with releases of hazardous substances (for example, HRS considers direct contact of people with contaminated soils, contamination of aquatic food chain, three broad types of human toxicity, potential for air contamination, sensitive environments that include wetlands, endangered species, and environments designated by various Federal and State agencies).
- HRS is the most carefully developed, thoroughly peer-reviewed, and widely applied of the targeting schemes to be reviewed under the waste minimization targeting effort. In the developmental stage of the HRS, components of the model were reviewed by the Science Advisory Board (SAB). Additionally, the final version of the HRS model is currently under review by the National Academy of Sciences. Carrying the weight of these reviews as well as public notice and comment, the HRS enjoys more regulatory and scientific credibility than the other schemes.
- HRS is considered to be a screening tool, not a detailed risk assessment, and is thus an
  appropriate targeting strategy.

- In 'full' form, HRS is relatively complex and data-intensive.
- HRS is site-oriented, not waste-oriented; the overall scores relate more to releases from sites and threats to nearby receptors. Some components of the HRS, however, may be useful to waste minimization targeting. For example, the waste characteristics factors, e.g., toxicity and persistence, can be used for scoring waste streams as generated.
- HRS is not specifically designed for generators of hazardous wastes.
- HRS is not designed to consider combustion or other treatment processes with respect to "quantitative release." An observed release has to be attributable to a source on the site in order to be scored; combustion and other active treatment units are not regarded as sources in the HRS.

Exhibit 1. Hazard Ranking System: Factors and Factor Categories

Pathway	Factor Categories		
- Talaway	Likelihood of Release	Waste Characteristics	Targets
Ground Water Migration Pathway	Observed Release or Potential to Release Containment Net Precipitation Depth to aquifer Travel time	Toxicity/Mobility Hazardous Waste Quantity	Nearest Well Population Resources Wellhead Protection Area

	Factor Categories		
Pathway	Likelihood of Release	Waste Characteristics	Turgets
Surface Water Migration Pathway	Overland Flood Component	Drinking Water Threat	Drinking Water Threat
	Observed Release	Toxicity/Persistence/Mobility	Nearest Intake
	or	Hazardous Waste Quantity	Population
	Potential to Release		Resources
	By Overland Flow:	Human Food Chain Threat	
	Containment		Human Food Chain Threat
	Runoff	Toxicity/Persistence/Mobility/	
	Distance to Surface Water	Bioaccumulation	Food Chain Individual
I		Hazardous Waste Quantity	Population
	By Flood:	_	
	Containment	Environmental Threat	Environmental Threat
	Flood Frequency		
		Ecosystem Toxicity/Mobility/	Sensitive Environments
	Ground Water to Surface Water	Persistence/Bioaccumulation	
	Component	Hazardous Waste Quantity	
	Observed Release		
	or		
	Potential to Release		
	Containment	·	
	Net Precipitation		
	Depth to Aquifer		
	Travel Time		

Exhibit I (continued). Hazard Ranking Bystem: Factors and Factor Categories.

Dash	Factor Categories		
Pathway	Likelihood of Release	Waste Characteristics	Targets
Soil Exposure Pathway	Resident Population Threat  Observed Contamination  Nearby Population Threat  Attractiveness/Accessibility Area of Contamination	Resident Population Threat  Toxicity Hazardous Waste Quantity  Nearby Population Threat  Toxicity Hazardous Waste Quantity	Resident Population Threat  Resident Individual Resident Population Workers Resources Terrestrial Sensitive Environments  Nearby Population Threat  Population Within 1 Mile Nearby Individual
Air Migration Pathway	Observed Release  or  Potential to Release  Gas  Gas Containment  Gas Source Type  Gas Migration Potential  Particulate  Particulate Containment  Particulate Source Type  Particulate Migration  Potential	Toxicity/Mobility Hazardous Waste Quantity	Resources Population Within 4-Mile Radius Nearest Individual Sensitive Environments

Exhibit 2. Targeting Criteria Used in the Hazard Ranking System.

TARGETING CRITERIA			
Direct Risks			
Waste volume	1		
Waste/constituent toxicity Human toxicity Ecological toxicity	11		
Constituent concentration or mass	1		
Waste type (e.g., solvents)			
Number of generators			
Waste management practices	1		
Releases to environmental media	1		
Potential for constituent transport	1		
Potential for cross-media transfer	1		
Potential for human exposure: residential settings occupational settings	1,		
Potential for ecological exposure	1		
Indirect or Acute Risks			
Ignitability, corrosivity, reactivity			
Ozone depletion potential	·.		
Global warming			
Other Targeting Criteria			
Hazardous waste management capacity			
Technical/administrative feasibility			
Permitting/enforcement factors			
Cost savings			
Other factors  Resource exposure (e.g., if aquifer being evaluated is used for drinking water or irrigation)	,		
Wellhead Protection Areas	1		

# Toxics Release Inventory Environmental Indicators Methodology<sup>12</sup>

# Purpose of the Methodology/System

- Proposed by EPA Office of Pollution Prevention and Toxics (OPPT) for use in tracking changes in human health and environmental risks posed by chemicals released to the environment.
- Will "allow EPA to measure its successes in implementing environmental protection and pollution prevention programs, and to formulate strategic plans for improving the course of future environmental progress."

#### System Description

- Calculates a national index of chronic health risks to human populations based on estimated exposures to environmental releases of TRI chemicals. Comparison of indices from year to year reveals whether risks are increasing or decreasing over time.
- Each facility-reported rele: of each TRI chemical to each environmental medium is weighted by toxicity, exposure potential, and the size of the exposed population to produce a risk-related "subindex." Basic steps:
  - 1. Determine location of TRI-reporting facility or facility that receives its wastes;
  - 2. Using a geographically-indexed database, match geographic and demographic features to facility location to derive site-specific environmental (e.g., stream velocity) and exposure (e.g., number of people using private drinking water wells) information for subsequent modeling;
  - 3. Estimate environmental concentrations at exposure points using TRI-reported release and generic and site-specific environmental and exposure information as input to mathematical models;
  - 4. From the exposure concentration, calculate dose using standard exposure assumptions;
  - 5. Assign exposure score based on calculated dose and degree of uncertainty;

<sup>12</sup> Documentation reviewed:

U.S. EPA, May 22, 1992, Toxics Release Inventory Environmental Indicators Methodology, Draft Report prepared for the Office of Pollution Prevention and Toxics by Abt Associates, Inc.

- 6. Multiply exposure score by size of exposed population and adjust for uncertainty; and
- 7. Multiply result by toxicity score reflecting chemical's potency and weight-ofevidence classification.
- Subindices for about 500,000 combinations of facility, chemical, and medium are summed to derive the national index.
- Five types of direct TRI releases (or "media") covered: air releases, surface water releases, on-site land releases, discharge to POTWs, and off-site transfers. The methodology for incorporating exposure into the indicator considers cross-media transfers also.
- Methods for estimating exposures to different types of releases vary significantly in level of rigor and use of generic vs. site-specific data.

#### System Evaluation

#### Scope of coverage:

- The TRI indicator includes all facility-reported releases of TRI chemicals (about 240-320 chemicals) to the environment, except those meeting certain criteria for exclusion.
- Based on certain criteria, some TRI chemicals proposed to be excluded from the TRI indicator (e.g., non-TSCA chemicals, chemicals with no reporting or zero reporting)

# Media/pathways addressed:

- Air releases:
  - -- Inhalation exposure
- Surface water releases:
  - -- Drinking water (surface water)
  - -- Fish ingestion
- Discharge to POTWs:
  - Efficent:
    - Drinking water (surface water)
    - Fish ingestion
  - Treatment:
    - Volatilization/inhalation
  - Sludge management:
    - Route depends on management practice
- On-site land releases:

- Drinking water (ground water)
- Volatilization/inhalation (actually reported under TRI air releases and handled along with other air releases)
- No release assumed if management is in a RCRA Subtitle C unit
- Off-site transfers:
  - -- Incineration
    - Inhalation
  - -- Land management
    - Volatilization/inhalation
    - Drinking water (ground water)
  - -- No release assumed if management is in a RCRA Subtitle C unit
- Cross-media transfers considered. For instance, on-site releases include releases to landfills, surface impoundments, land treatment units and underground injection. To evaluate exposure from these releases, the indicator methodology models "cross-media" transfers also, e.g., leaching from landfill to groundwater and volatilization. For releases to POTWs, the methodology models volatilization and adsorbtion to and subsequent volatilization from sludge.

# Types of targeting criteria used (see Exhibit 1):

- Overall measure: chronic health risks to human populations (not to individuals)
- Incorporates standard risk assessment factors and methods

# Data requirements:

- Constituent toxicity (all endpoints specified in SARA Section 313): from IRIS; HEAST; structure/activity relationships and other methods
- Releases to environmental media: from TRI database
- Estimation of environmental concentrations and size of exposed population:
  - Site-specific data needs and sources:
    - Facility location from TRI
    - Geographically-indexed environmental and exposure information from BGREACH file, a SAS file developed for this project and housed on the EPA mainframe
    - RCRA regulatory status from TRI database, RCRIS
  - Generic inputs/default values derived using professional judgment and information from other EPA studies

# Applicability to waste minimization targeting:

- System is designed to be used in tracking successes in pollution prevention and therefore seems appropriate for use in waste minimization/pollution prevention targeting.
- Cross-media transfers are covered.
- Subindices can be calculated to evaluate contribution to national index of individual chemicals, regions, states, industries, or release pathways.
- Method is based on standard risk assessment paradigm and is relatively rigorous.

- Very difficult to use system to track only hazardous waste-related releases and risks without major programming changes:
  - -- Waste management in Subtitle C units is excluded (except for on-site incineration in Subtitle C units, which is handled along with other on-site air releases and may be difficult to separate out).
  - -- Unclear whether incinerator ash management is covered at all.
- Methods for calculating exposures from land-based waste management are relatively unsophisticated and use mostly generic inputs (waste concentration calculated by dividing TRI-reported releases by industry-wide average annual waste generation; leachate concentrations using a partitioning approach; ground-water concentrations using a generic dilution and attenuation factor (DAF).
- Ecological risks not covered, but are proposed to be added.

Exhibit 1. Targeting Criteria Used in the TRI Environmental Indicators Methodology.

TARGETING CRITERIA	
Direct Risks	
Waste volume	
Waste/constituent toxicity Human toxicity Ecological toxicity	1,
Constituent concentration or mass	1
Waste type (e.g., solvents)	
Number of generators	
Waste management practices	1
Releases to environmental media	1
Potential for constituent transport	1
Potential for cross-media transfer	
Potential for human exposure: residential settings occupational settings	1
Potential for ecological exposure	
Indirect or Acute Risks	
Ignitability, corrosivity, reactivity	
Ozone depletion potential	
Global warming	. # "
Other Targeting Criteria	
Hazardous waste management capacity	
Technical/administrative feasibility	
Permitting/enforcement factors	
Cost savings	
Other factors	

# Toxics Release Inventory (TRI) Risk Screening Guide 13

#### Purpose of the Methodology/System

- TRI a national computerized data base containing data submitted by industry on the hazardous chemicals manufactured, used, stored, processed, or released to air, surface water, POTWs, or land.
- The TRI risk screening guide (system) is a framework for initial, screening-level analyses to identify, from among all TRI submissions for a particular geographical area or community, those risk scenarios, facilities, or chemicals for follow-up investigation.

#### System Description

- Qualitative or relative expression of risk (i.e., high, moderate, or low) derived by evaluating various chemical-specific and site-specific factors. These factors, listed in Exhibit 1, are evaluated in three different components that comprise the system:
  - Toxicological potency. Incorporates measures of both the nature of the adverse human health or ecological effects (hazard identification) and the magnitude of these effects at specific exposure levels (dose-response relationships). Uses readily available EPA estimates of toxicological potency, which include reportable quantities (RQs), threshold planning quantities (TPQs), cancer potency factors, reference doses (RfDs), and ambient water quality criteria (AWQC).
  - Exposure evaluation. Qualitatively evaluates data for two key aspects of exposure plausible exposure pathways and potential environmental levels based on site-specific and chemical-specific factors. The potential environmental levels refer to the characterization of amounts and concentrations of a chemical in the environment at points of contact with populations of interest.
  - Risk characterization. Qualitatively combines the toxicological potency assessment and the exposure evaluation (i.e., in terms of relative probability of harm such as high, moderate, or trivial) to identify facilities, populations, and chemicals that warrant further investigation. Produces a profile of scores for each medium.

<sup>13</sup> Documentation reviewed:

U.S. EPA. 1989. Toxic Chemical Release Inventory Risk Screening Guide (Version 1.0), Volume I: The Process. Prepared by the Office of Toxic Substances, U.S. Environmental Protection Agency, Washington, D.C., EPA 560/2-89-002.

Exhibit 1: Factors Used in TRI Risk Screening

Toxicological Potency	Human Health Effects	Ecological Effects	.9	Dose-response Relationships	7
	Curchogenicity     Heritable gene and chromosome mutations     Neurotonicity     Remoducibekteetkumental terricity	Environmental toucity     Toucity and persistenc     Toucity and beoseeum	Environmental toxicity Toxicity and persistence Toxicity and besaccumulation Other sombrant effects	Properties of individual chemicals     Duse, frequency, & duration of exposure     Route of expusure	
	Other chronic effects Adverse acute effects			Individually susceptibility	
Exposure Evaluation	Pleasible Exposure Pathways		Potential Environmental Levels	nental Loveis	
	Site-specific:		site-specific:	÷	
	Location of release     Population(s) exposed     -human or ecological		Geographic d     Physical trans	Geographic distance to populations Physical transport characteristics	
	-scrstive subcoculations		Chemical specific:		
	Uses of contaminated media		Physical fate:	Physical fate and transport characteristics	
	Cicographic distance to populations     Physical transport characteristics		Emvironmental     Rute of release	Finvironmental transformation characteristics Rate of release	
	Chemical resertion		<ul> <li>Quantity of release</li> </ul>	kasc	

	L	
Risk Characterization	•	Qualitative combination of toxicological potency and emorate evaluation
Component	•	Includes characterizing the nature of uncertainties and quality of data

Physical fate and transport characteristics Environmental transformation characteristics

Chemical-specific

#### System Evaluation

#### Scope of coverage:

- Constituents: Covers over 300 chemicals and/or chemical categories (i.e., components of mixtures) that are subject to TRI reporting.
- Releases: Covers releases from facilities:
  - -- classified in Standard Industrial Classification (SIC) codes 20 through 39; and
  - -- that employ the equivalent of 10 or more full-time individuals; and
  - that manufacture (including import) or process any of the Section 313 chemicals or chemical categories in amounts greater than 75,000 pounds in 1987; 50,000 pounds in 1988; or 25,000 pounds in 1989 and subsequent years; or who use any listed chemical or chemical category in any other way (other than manufacture) in amounts greater than 10,000 pounds in 1987 and subsequent years, including processing or importing the listed chemical or chemical category.
  - -- Only routine releases (i.e., releases occurring during normal industrial operation) are addressed: "burst" or accidental, rapid releases are not addressed.

#### Media addressed:

- All releases to:
  - -- air (fugitive and point-source)
  - -- land (injection, landfills, surface impounding, landspreading)
  - .. surface water and publicly owned treatment works (POTW)

#### Types of targeting criteria used:

- Targeting based on human health and ecological risks, which in turn are assessed based on site- and chemical-specific factors listed in Exhibit 1.
- Local, State, and Federal laws, public concern, control technologies, economics, and politics are additional criteria that can be used in the risk-management phase to fully characterize risks from routine TRI releases.
- Exhibit 2 summarizes the key targeting criteria used in this system.

#### Data requirements:

• Data needed for the screening are less detailed than data required for a formal risk assessment.

- Relevant data are obtained from the TRI data base and from appendices to the risk screening guide.
- Supplementary information is required to fully characterize risks for TRI releases and to
  put these risks into perspective with those from other chemical releases in the community.

# Applicability to Waste Minimization Targeting:

#### Pros

- System is a screening-level tool that conforms with the risk assessment paradigm.
- System documentation provides toxicity indices for TRI chemicals and methods to interpret and use TRI releases data.

- System is designed around TRI data, which does not cover all chemicals, wastestreams, or industrial sectors. For example, not all industrial releases of listed chemicals are covered by the TRI reporting requirements (facilities with fewer than 10 employees, industries outside the specified SIC codes, and industries using less than the threshold quantities).
- Release data is difficult to use for exposure modeling. Releases reported by industry are summary data reported in pounds per year; no specific information on frequency, duration, concentration, or peak release is required. Reporting form does not specify location of point sources or identity of chemicals comprising mixtures and compound classes.

Exhibit 2: Targeting Criteria Used in the TRI Risk Screening Guide

TARGETING CRITERIA	
Direct Risks	
Waste volume (pounds)	1
Waste/constituent toxicity Human toxicity Ecological toxicity	<b>*</b>
Constituent concentration	
Waste type	
Number of generators	
Waste management practices	
Releases to environmental media	1
Potential for constituent transport	1
Potential for cross-media transfer	
Potential for human exposure: environmental settings occupational settings	**
Potential for ecological exposure	
Indirect or Acute Risks	
Ignitability, corrosivity, reactivity	
Ozone depletion potential	
Global warming	-
Other Targeting Criteria	
Hazardous waste management capacity	
Technical/administrative feasibility	
Permitting/enforcement factors	1
Cost savings	
Other factors	

# APPENDIX 8 HRS HAZARD DATA AND PATHWAY SCORES FROM SUPERFUND CHEMICAL DATA MATRIX

ž

		, ,				HRS Pathway Sc	ores				
Chomical	CAS No	Toxicity	OW Mob	Air Mob	Persist	PCBioacc	EnvBioacc	Ecotox	OW	Air	SW/OP/DWT
		•							(Tox/Mob)	(Tox/Mob)	(Ton/Per)
1.1.1.2-tetrachloroethane	630-20-6	1.00E+02	1 00E-02	1 00E+00	4 00E-01	5 00E+01	5 00E+01	1 00E+01	1 00E+00	1 00E+02	4.00B+01
1,1,1-trichloroethane	71-55-6	1 00E+01	1.00E-02	1.00E+00	4 00E-01	5.00E+00	5 00E+00	1 00E+01	1 00E-01	1.008+01	4.00E+00
1.1.2.2-tetrachloroethans	79-34-5	1 00E+01	1.006-02	1.00E+00	4.00E-01	5.00E+00	5 00E+00	1 00E+02	1 00E 01	1 005+01	4.00B+00
1.1.2-trichloreethens	79-00-5	1 00E+03	1.00E-02	1.00E+00	4.00E-01	5.00E+01	5 00E+01	1.00E+01	1 001:+01	1.00£+03	4.00B+02
1.1.2-trichlore-1,2,2-trifluoroethane	76-13-1	1.00E+00	1 00E-02	1 005+00	4.00E-01	5 00E+01	5.00E+01	NV	1 001: -02	1.00B+00	4 00B-01
1.1-dichloroethane	75-34-3	1.005+01	1.00E+00	1.006+00	4.00&-01	5.00E+00	5 006+00	NV	1.00E+01	1.00E+01	4.00B+00
I .I - dichloroethylene	75-35-4	NV	NV	NV	NV	NV	NV	NV	NV	NV	NV
1,2,3-trichloropropase	96-19-4	1.005+02	1.00E-02	1.00E+00	4.00E-01	5.00E+00	5 00E+00	1 00E+01	1 001:+00	1 00E+02	4 00B+01
1,2,4,5-tetrachlorobenzeno	95-94-3	1.00E+04	1.00E-04	2.00E-01	1.00E+00	5.00E+03	5.00E+03	1.00E+02	1.001:+00	2.00E+03	1.00E+04
1,2,4-trichlorobenzene	120-82-1	1.006+02	1.00E-02	1.005+00	4.00E-01	5.00E+02	5.00E+02	1.00E+03	1 00E+00	1.00E+02	4.00B+01
1.2-dibromoethass	106-93-4	1.00E+04	1.005+00	1.00£+00	4.00E-01	5 00E+00	5 00E+00	NV	1 00E+04	1 00E+04	4.00B+03
1,2-dibresso-3-chleropropens	96-12-8	1.005+04	1.006-02	1.006+00	1.00E+00	5.00E+01	5 00E+01	NV	1 00E+02	1.00E+04	1.00B+04
1.2-dichlorobenzone	95-50-1	1.00B+01	1 00E-02	1.00E+00	4.00E-01	5.00E+01	5.00E+01	1.00E+02	1.00E-01	1.00E+01	4.00E+00
1.2-dichloroethons	107-06-2	1.00B+02	1.00E+00	1.005+00	4.00E-01	5.00E+00	5.00E+00	1.00E+00	1.00E+02	1.005+02	4.00E+01
1.2-dichloroethylens	156-60-5	NV	NV	NV	NV	NV	NV	NV	NV	NV	NV
1.2-dichloropropane	78-87-5	1.00E+03	1.00E+00	1.00E+00	4.60E-01	5.00E+01	5.00E+01	1.00E+01	1.00E+03	1.00E+03	4 008+02
1.3-dichlorobeaseas	541-73-1	NV.	1.00E-02	1.00E+00	4.005-01	5.00E+01	5.00E+01	1.00E+02	NV	NV	NV
1,3-dichloropropens	542-75-6	1.00E+04	1.005+00	1.00E+00	4.006-01	5.00E+00	5.00E+00	1.00E+03	1 00E+04	1.00E+04	4.00E+03
1.4-dichlorobenzens	106-46-7	1 00E+01	1 00E-02	1.00E+00	4.00E-01	5.00E+01	5.00E+01	1.00E+02	1 00E-01	1.006+01	4.00E+00
2,3,4,6-tetrachlorophenol	58-90-2	1:00E+02	1.006-04	2.00E-01	1.00E+00	5.00E+02	5.00E+03	1.00E+03	1.00E-02	2.005+01	1 00E+02
2.d.5-TP (Silven)	93-72-1	1.006+02	1.00E-02	2.00E-03	4.00E-01	5.00E+02	5.006+02	1.00E+04	NV	2.00E-01	4.008-01
2,4,5-trichlorophenol	95-95-4	1.00E+01	1.00E-02	2.00E-01	1.00E+00	5.00E+02	5.00E+03	1.00E+03	1.00E-01	2.00E+00	1.00E+01
2,4,6-trichlorophenol	88-06-2	1.00E+01	1.00E+00	2.00E-01	1.00E+00	5.00E+02	5.00E+04	1.00E+03	1.00E+01	2.00E+00	1.00E+01
2,4-D	94-75-7	1.00fi+02	1 00E-02	2.00E-03	1.005+00	5.00E+01	5.00E+01	1.00E+02	NV	2.00E-01	1 00E+02
2,4-dichlorophenol	120-83-2	1.00B+03	1.00E-02	2.00E-01	1.00E+00	5.00E+02	5 00E+02	1.00E+02	1 00E+01	2.00E+02	1 00E+0
2-chloremethylexirans	106-89-8	1:00E+04	1.00E+00	1.00E+00	4.00E-01	5.00E-01	5.00E-01	1.00E+01	1.00E+04	1.00E+04	4.00E+03
2-chlorenaphthalens	91-58-7	1.00E+01	1:00E-04	1.10E+01	1.00E+00	5.00E+02	5.00E+02	NV	1.00E-03	1 10E+02	1 00E+01
2-chlorophanel	95-57-8	1.00E+02	1.00B-02	1.00B+00	4.00E-01	5.00E+02	5.00E+02	1.00E+02	NV	1.00E+02	4.00B+01
3.3-dichlorobentidika	91-94-1	1.00B+02	1.00E-04	2.00E-04	1.00E+00	5.00E+02	5.00E+02	NV	1.00E-02	2 00E-02	1 00E+02
acenaphtheae	83-32-9	1.00E+01	1:00E-02	2.00E-01	4.00E-01	5.00E+02	5 00E+02	1.00E+04	1.00E-01	2.00E+00	4 006+02
scannehthylone	208-96-8	NV	1.00B-02	2.005-02	1.00E+00	5.00E+02	5.00li+02	NV	NV	2.00E-00	NV
acetaldohyda	75-07-0	1.00E+03	1.00E+00	1.00E+00	4.006-01	5.00E-01	5.008-01	1.00E+01	1.00E+03	1 00E+03	4 00E+02
acetone	67-64-1	1.006+01	1 00E+00	1.00E+00	4.00E-01	5.00E-01	5.00E-01	1.00E+02	1.00E+01	1.00E+01	1
acotonitrile	75-05-8	1.005+02	1.00E+00	1.00E+00	1.006+00	5.00E-01	5.00E-01	4 00E+00	1 001:+02	1 00E+02	4 00E+00
acetophenone	98-84-2	1.00E+01	1.00E+00	1.00E+00	4.00E-01	5.00E+00	5 00E+00	1.00E+00	1.00E+01	1.00E+02	1.00E+02
scroleia	107-02-8	1.00E+04	1.00E+00	1.00E+00	4.00E-01	5 00E+02	5 00E+02	1.00E+04	1.00E+01	1.00E+01	4 00E+00
acrylamide	79-06-1	1.00E+03	1.00E+00	2.00E-01	4.00E-01	5.00E+00	5.00E+00	1.00E+01			4 00E+03
acrylic acid	79-10-7	1.00E+03	1.00E+00	1.00E+00	1.00E+00	5.00E-01	5.00E-01	NV	1.00E+03 1.00E+03	2.00E+02 1.00E+03	4 00E+02
acrylonitrik	107-13-1	1.00E+03	1.00E+00	1 00E+00	4.00E-01	5.00B+01	5.00E+01	1.00E+02	1.00E+03		1 00E+03
allyl chloride	107-05-1	NV	NV	NV	NV	NV	NV NV	NV	NV	1.00E+03	4 00E+02
alpha-BHC	319-84-6	NV	NV	NV	NV	NV	NV	NV	NV NV	NV NV	NV
emmonia	7664-41-7	1 00E+02	1 00E+00	1 00E+00	7.00E-04	5.00E-01	5 00E-01	1 00E+01	1.00E+02		NV 7 COLUMN
aniline	62-53-3	1 001:+04	1.00E+00	1 00E+00	1.00E+00	5 00E+00	5.00E+02	1.00E+01		1 00E+02	7 0015-02
anthracene	120-12-7	1.00E+01	1.00E-04	2.00E-03	4.00E-01	5.00E+03	5.00E+03	1 00E+04	1 00E+04 1 00E-03	1,00E+04 2 00E-02	1 00b+04 4 00B+00

# finhibit 1. HES Hazard Data and Pathway Scores for Halogenated Organics

						HRS Pathway Scores	COCCO				
Occursed	CAS No	Toronto	A W	Aii Mak		50					
		- outry	7	211 9100	£ CI MONE	L Browc	Emviliance	Ecolox	WD	) I	SW/OP/DWT
arocler (016	12474-11-2	NV.	NO.	AIN	No.				(Ton/Mub)	(Jor/Mop)	(TowPas)
aroclor #221	11104-28-2	£ :	2 7		Į .	2	2	N	Z	ž	NV
armcler 3232		¥ ;	<b>.</b>	2 2	2 4	2	×	N	Z	×	74
artist 1248	111111111111111111111111111111111111111	¥ 7		74	2	N	×	Z	Z	Y Y	Z
174 m	10-21	2	200	Z	2	z	VV	N	2	ž	Z
analy 1740	3	2	N	Z	N N	×	N	7	N	N.	Z
Bracks 1480	11000 12-5	N	N	N	Z	N.	N	N	z	¥	<b>Z</b> <
	¥ 55-3	1 005+03	1 00E-04	2.00E-04	1.00E+00	5.00E+04	5.00E+04	1 001:+04	1.001:-01	2.00E-01	1.00E+03
	71-43-2	1.0011-02	1.00E+00	1.00E+00	4.00E-01	5.00E+03	5.00E+02	1 005+03	1 001:-02	1.00E+02	4.00E+01
wanted in Property and	30-32-0	1.008+04	1.006-04	2.00E-04	1.00E+00	5.00E+04	5 00l5+04	1.005+04	1.001:-00	2.00E+00	I ODE-OA
Best Culturations	207-04-9	Z	1.00E-04	2.00E-04	1.005+00	5.00E+04	5 005+04	N	Z C	N	200
Bearing Caterions	100-44-7	1.006+02	1.00E-02	1.00E+00	7.00H-02	5.00E+01	5.006+01	1.00E+02	Z	i Milan	7005.00
bota-BHC	319-85-7	ş	N.	Z C	Z	Z	z	Z	2 :	NV.	medim.
bis (A-chloroshoxyi) methane	111-91-7	1.00E+02	1.00E+00	6.00E+00	1.00E+00	5.00E-01	5.00E-01	Z	1 main	\$ 000 Lave	- Ne
the (2-staythery)) philastate	117-81-7	1.00E+02	1 00E-04	2.00E-03	1.005+00	5.00E+02	5.00E+04	1.00E+03	1 001:-02	2005-01	1.000
DE CALCADOR DE CAL	75-27-4	1.00E-02	1.00E-02	1.00E+00	4.00E-01	3.00E+01	3.00H-01	z	NV	1 001:-02	70-000
	73-23-2	Z	N	Y	Z	٧ĸ	Z C	Z	z	N .	NV.
carbon disulfida	74-83-9	1 001:+03	1.00E-02	1.00E+00	4.00E-01	5.00E+00	5 00H+00	1.001:+04	100F+01	1.00E+03	4 00E+02
carbon tetrachloride	4 2 2 4	1 005-03	20-800.1	1.005+00	4.00E-01	3.00E+01	5.00E+01	1.00E+02	1.001-01	1.00E+03	4 00E+02
chioni	75-27-4	1 005-03	1 000-02	1.000	4.00E-01	5.0011-01	5.00E+01	1.00E+02	1.00E-01	1.00E+03	4 00E+02
chlordage	57-74-0	1 00000	1.00H	1.1015-01	1.00E+00	5.00E+00	5.00E+00	×	1.00E+03	1.106+04	1 006-01
chlorobeazene	101-01-7	1.000	00E-00		1.00E+00	5.00E+04	5.00E+04	1.00E+04	Z C	6.00E+04	1.00H-04
chlorosthanc	75-88-3	2000	1.005-02		7.001:-04	5.00E+01	10+300 \$	1.00E+03	N C	1.70E+03	7 0015-02
c Morodorm	27-14-1	1.005-00		West of the second	7.00b-04	5.00E+00	5.00E+00	24	1.005+00	1.001-00	7 00E-04
chloresethano	74-17-1	20000	1.005-00	10-807	4.00E-01	5.00E+00	5.006+00	1.00E+01	1.00E+02	1.705+03	4 00 E+01
chiorewothyl methyl other	107-45-5	1.000.01		10-250	7.00b-04	5.00E+00	5.00B+00	1.00E+00	1.00H-01	1.705+02	7.00E-01
creed (m-creed) •		1 00R-01	00000	10.00	/.OUb-04	3.00E-01	5.00E-01	2<	1.00E+03	1.70E+04	7.008-01
сгужена	218-01-0	Y	100 D		1.005+00	5.00E-00	3 00E+00	1.00E+02	1.005-01	1.001-01	005-01
Cumany	3		1.000	7.00H-04	1.00E+00	5.00H-02	5.00E-03	1 00E+03	Z <	Z	
cyanoges chleride	5	TO TO TO	70-1300.1	1.00t±00	4.00E-01	5.00E+02	5.00E+02	1.00E+02	1.006+01	1.00E+03	4 Mileson
cyclehoxanona		2	NO	24	Z	2<	N C	N K	Z C	Z	TOUR P
DDD	77-4-	00400	1.000	1.00H+00	1.00E+00	5.00E+00	5.00E+00	1.00E+00	1 00E+00	1.00E+00	35.70
DDT	5	1.005.01	200	2.00E-03	1.00E+00	5 00E+04	5.00E+04	1 0015-04	1 006-02	2.00E-01	006-00
dichloradiffuoromethase	75-71-1	1.000.00		2.00E-03	1.00E+00	5.00E+04	5 00E+04	1 005+04	1 00E-01	2.00E+00	20-200
dichlerockyl ether		100000	70-900.1	1.00±±00	4.00E-01	3.00E+01	5.00E+01	N C	1.00E-01	1008+01	4 005-00
dichloratoprapy) ether		2	2	Z	Z	z <	N C	<b>X</b>	Z K	Z .	- COLOR 100 COLO
dichloromothy! other		2	Z	×	z	N	z	X <	Z :	Z :	2 2
dieldrig	342-84-1	NV	Z	7	Z	N V	Z <	Z <	Z :	2 2	2
distant mirror	60-57-1	1 00E+04	1.00E-04	2.00E-03	1.005+00	5.00E+04	5.00E+04	1 005-04	2 2	N	Z
and a second sec	86-30-6	1.00E+01	1.00E-02	2.00E-02	1.00E+00	5 00E+02	\$ 00E-03	1.001.00	7	2 00E+01	1 001:-04
-presentation y GOLD	106-89-8	7	Z Z	z <	Z	NV.	7002007	70+300.1	10-400	2 00E-01	1 00E+01
cmyl accuse	141-78-6	1 00E+00	1.001:+00	00:-00	4 005-04	A NO.	2	Z	×	ž	7
ciayi ocalcae	100-41-4	1.00E+01	1 001:-02	-00	100E-01	\$ 000.01	10-300 c	1 00:+00	1 001:-00	1 00E+00	4 00E-01
cinyien: aibromide	106 93-4	z	z	Z	N.V.	3.000	10+300.0	1 00E+02	10-3001	1.001-01	4 00h-00
cunyica: glycol	107-21-1	00+300 I	3	2000	-	24	2	Z	Z Z	z -	1
					355	300				• • • •	z <

	1					HRS Pathway Sc	ores				
homical	CACNE	Toxicity	OW Mob	Air Mob	Persist	FCBioacc	Linvilionece	Exotox	- ow	Air	SW/OF/DW
	CAS No	toricity	G.V MIOU	2100	. 3				(Tox/Mob)	(fox/Mob)	(Tox/Pc
41.4	110 80-5	1 001:+01	1 00E+00	1 00E+00	1 001:+00	5.00E-01	5 00E-01	1.00E+00	1.00E+01	1.001:+01	1 00B+0
thylene glycol monomethyl ether	60-29-7	1 00E+01	1 001:+00	1 001:+00	4.00E;-01	5 00E-01	5 001:-01	1 00E+00	1 00E-01	1.00E+01	4 00B+0
thyl other	204-44-0	1 00E+02	1 00E-04	2 001:-04	1 00E+00	5 00E+03	5 00E+03	1 00E+04	1 00E-02	2 00E-02	1.00E+
Nuoranthene	24-73-7	1 00E+02	1 00E-02	2 00E-01	1 0015+00	5 00E+03	5 00E+03	1 00E+03	NV	2 00B+01	1.00B+
Nuoreae	50-00-0	1 00E+01	1.00E+00	1 60E+00	1.00E+00	5 00E-01	5 00E-01	1 00E+02	1 00E+01	1.00E+01	1.00₽↔
formaldchydc	76-44-8	1 00E+03	2.00E-02	2.00E-02	1.00E+00	5 00E+03	5 00E+04	1 00E+04	2 001:+01	2.00B+01	1.00B+
septachlor	1024 - 57-3	1 00E+04	1 00E+00	2 00E -02	1.00E+00	5.00E+00	5 00E+04	1 00E+04	1 00E+04	2 00E+02	1 00E+
acptachlor epoxide	118-74-1	1.00E+03	1.00E-04	2.00E-02	1 00E+00	5 00E+03	5.00E+04	1.00E+01	1.00E-01	2 00E+01	1.00E+
hexachlorobenzene	87-68-3	1 00E+03	1.00E-04	2.00E-01	1 00E+00	5 00E+01	5 00E+03	1.00E+04	1 00E-01	2 00E+02	1.00B+
hexachlorobutadiene (hexachloro-1,3-butadeene)	77-47-4	1.00E+04	1 00fi-02	2 00E-01	1 00E+00	5 006+03	5 00E+01	1 00E+04	1 00E+02	2.00E+03	1.00B+
he xachlorocyclopeniadions	67-72-1	1.00E+03	1.00E-02	1.00E+00	4.00E-01	5.00E+02	5.00E+02	1.00E+03	1.00E+01	1.00E+03	4 00B+
he xac hioroethans	77-47-4	1.00E+04	1 00E-02	2.00E-01	1.00E+00	5 00E+03	5.00E+01	1 00E+04	1.00E+02	2 00E+03	1 00E+
hexachloropentadiene	7647-01-0	1.00E+03	1.00E+00	NV	4 00E-01	5 00E-01	\$ 00E-01	1 00E+00	1 00E+03	NV	4 00E+
hydrochloric scid	7783-06-4	1.00E+04	1.00E+00	1.00E+00	4.00E-01	5 00E-01	5 00E-01	1.00E+03	1 001:+04	1 00E+04	4 00E+
hydrogen sulfide	78-83-1	1 0015+01	1.00E+00	1.00E+00	4.00E-01	5.00E-01	5.00E-01	1.00E+01	1 00E+01	1 00E+01	4 00É+
sobutanol	58-89-9	1 00fi+04	1 00E-03	2.00E-02	1.00E+00	5.00E+02	5.00E+02	1 00E+04	1.00E+01	2.001:+02	1 008
lindene	67-56-1	1 00E+00	1.00E+00	1 00E+00	1.00E+00	5 00E-01	5.00E-01	1.00E+00	1.005+00	1.00E+00	1.00E
mothanol	72 43 5	1 00E+02	1.00E-04	2.00E-03	1.00E+00	5.00E+04	5.00E+04	1.00E+04	1.00E-02	2 00E-01	1 00E
methoxychlor	79 22 1	1 00E+02	NV	1.00E+00	4.00E-01	5.00E-01	5 00E-01	NV	NV	1 00E+02	4.00E
methyl chlorocarbonate	101 14 4	1 001:403	1 00E-04	2.00E-04	4 00E-01	5.00E+02	5.00E+02	NV	1.00E-01	2 00E-01	4 00E
methylene-bis(2-chloroaniline), 4,4	75 09-2	1 00E+01	1 00E+00	1.00E+00	4.00E-01	5 00E+00	5,00E+00	1.00E+00	1.00E+01	1 00E+01	4 005
methylene chloride	78-93-3	1 00E+01	1 00E+00	1.00E+00	4.00E-01	5.00E-01	5.00E-01	1 00E+00	1.00E+01	1 00E+01	4 00E
methyl ethyl ketone	108-10-1	1.00E+01	1.00E+00	1.00E+00	4.00E-01	5 00E+00	5.00E+00	1.00E+00	1 00E+01	1 00E+01	4 00E
methyl subsityl kelone	80-62-6	1.00E+01	1 006+00	1.00E+00	4 00E-01	5.00E+00	5.00E+00	1.00E+00	1.00E+01	1.00E+01	4 00E
methyl methocrylate	71-36-3	1.00E+01	1.00E-02	1.00E+00	1.008+00	5 00E+00	5.00E+00	1.00E+00	1.00E-01	1006+01	1 005
a bitanol :	91-20-3	1.00E+02	1.00E-02	2.00E-01	4.00E-01	5.00E+02	5.00E+02	1.00E+03	1.00E+00	2 00E+01	4 00E
napidhelene	98-95-3	1.00E+03	1.00E-02	1.00E+00	1.00E+00	5.00E+00	5.00E+00	5.00E+00	1.00E+01	1	1 006
nstrubenzene pontachlorobenzeno	608-93-5	1.00E+03	1.006-02	2.00E-01	1.005+00	5.00E+03	5.00E+03	1.006+02	1.005+01	2.00E+02	1.00B
postachloroothass	76-01-7	1.00E+01	1.006-02	1.00E+00	4.00E-01	5.00E+01	5:00E+01	1.00E+02	1.00E-01	1 00E+01	4.00B
postackiorositrobestess	82-68-8	1.00B+03	1.00E-02	2.00E-02	1.00E+00	5.00E+02	5.00E+02	NV	1.00E+01	2.00E+01	1.00E
pontachiorophonol	87-86-5	1.006+02	1.005-02	2.00B-02	1.00E+00	5.00E+02	5.00E+02	1.00E+02	NV	2.00E+00	1.00B
p-Chlorosnitins	106-47-8	1.00E+03	1.00E+00	1.105+01	1.00E+00	5.00E+00	5.00E+00	1.00E+04	1.00E+03	1.10E+04	1.00B
phonanthrone	85-01-8	NV	1.005-04	2.00E-02	4.00E-01	5.00E+01	5.00E+03	1.00E+03	NV	NV	,
specol	108-95-2	1 00E+00	1.00E+00	1.00E+00	1.00E+00	5.00E+00	5.00E+00	1.00E+04	1.00E+00	1.00E+00	- 1.00B
pyreas	129-00-0	1.00E+02	1.00E-04	2.00E-03	1.00E+00	5 00E+01	5.00E+01	NV	1.00E-02		1 00E
pyridiac	110-86-1	1.00E+03	1.005+00	1.00E+00	1.00E+00	5.00E-01	5.00E-01	1 00E+02	1.00E+03	1 00E+03	1.00B
styreas	100-42-5	1.00E+01	1 00E-02	1.00E+00	4.00B-01	5.00E+01	5.00E+01	1.00E+02	1 00E-01	1.00E+01	4.00B
mulfuric acid	7664-93-9	1.00E+03	1 00E+00	NV	4.00E-01	5.00E-01	5.00E-01	1.00E+01	1 00E+03		4 00B
tolucas	108-88-3	1.00E+01	1.006-02	1.00E+00	4.00E-01	5.00E+01	5.00E+01	1.00E+02	1 00E-01		4 00E
toxaphene	8001-35-2	1 00E+03	1 00E-02	2 00E-03	1.00B+00	5.00E+04	5:00E+04	1 00E+04	1.00E+01		1.008
tetrachloroetheae	127-18-4	1.00E+02	1 00E-02	1 00E+00	4.00B-01	5.00E+01	5.00E+01	1.00E+02	1.00E+00		4 006
icirahydrofuran	109-99-9	1 001:+00	1 00E+00	1 001:+00	1.00E+00	5.00E+02	5 00E+02	NV	1.00E+00		1 008
Irans-1,2-dichloroethylene	156-60-5	1 00E+02	1 00E+00	1 001:+00	4.00B-01	5.00fi+01	5.00E+01	1 00E+00	1 00E+02	1	4 008
trichlorocthylene	79 01-6	1 001:+01	1.00E-02	1.00E+00	4 00E-01	5 00E+01	5.00E+01	1.00E+02	1 00E-01	1 00E+01	4 008

fixhibit 1 HRS Hazard Data and Pathway Scores for Halogensted Organics

Chemical					1	HRS Pathway Sco	orce				
	CAS No	Toucity	ów Mob	Air Mob	Persist	PCBioacc	LuvBioacc	Ecotox	GW	Vit	SW/OF/DW
Inchloromonofluorumethane	75 69 4	NV	NV	NV	NV I					(fox/Mob)	(Tex/Per
viayl chlorido	75-01-4	1 001:+04		- 1		NV	NV	NV	NV	אי	'И
nylene (m-nylene) **			1 00E-02	1 00E+00	7.00E-03	5 00E+00	5 00£+00	NV	1 001:+01	1.00E+04	7.00B+0
• • • •	106-36-3	1 00E+00	1 00E-02	1 00E+00	4.00E-01	5 00E+02	5 00E+02	1 00E+02	1 00E-02	1 00H+00	4.00B-0
xylcac (p-xylcac) **	106-42-3	1 006+01	1 005-02	1 00E+00	4 005-01	5 00E+01	4 00E+01	1 00E+02	1 00E-01	1 00B-01	4.00E+0
2,4-Directhyl phonol	0-0-0	1 00E+02	1 00E-02	2 00E-01	1 00E+00	5 00E+02	5 001i+02	1 005+02	1 00E+00	2 00E+01	1,00B-0
Benzo(n)anthracene	0-0-0	1.00E+03	1 00E+00	2.00E-04	1.00E+00	5.00E+04	5 00h+04	1 00E+04	1 00E+03	2 00E-01	1.00B+0
Butyl benzyl phthalate	0-0-0	1 00E+01	1 005-04	2.00E-03	1.00€+00	5 00E+02	5 001:02	1.00E+02	1 00E-03	2.00E-02	
Dibonz(a,h) anthracene	0-0-0	1 00E+04	1.006-04	NV	1.006+00	5.00E+04	5.00E+04	NV	1.00E+00		1.00B+0
Dimothyl phthalate	0-0-0	1 00E+00	1.00E-02	2 00E-01	1.00E+00	5.00E+01	5 00h+01	*** [		NV	1.00B+04
Pluorine	0-0-0	1 005+01	1.00E+00	1 006+00	4.00E-01	5 00E-01		1.00E+01	1.00E-02	2.00E-01	1.008+00
Maleic anhydride	0-0-0	1.00E+01	NV	1.00E+00	7.00E-04		5.00E-01	NV	1.00E+01	1 006+01	4.00B+00
Phornia	0-0-0	1.00E+04	1.00E-02			5 00E-01	5.00E-01	1.00E+00	NV	1 00E+01	7.00B-03
Phthalic anhydrade				2.00E-02	1.00E+00	\$.00E+03	5.00E+01	1.00E+04	1.00E+02	2.00E+02	1.006+04
Teluana dissocyanate	0-0-0	1.00E+00	1.00E+00	2.00E-02	4.00E-01	5 0 1 01	5.00E-01	NV	1 00E+00	2.00E-02	4.00E-0
	0-0-0	1.00E+03	иv	2.00E-01	4.00E-01	5.00E-01	5.00E-01	1.00E+00	NV	2 00E+02	4.00E+0
2,3,7,8-Totrachlorodibeazo(p)dioxin	0-0-0	1.00E+04	1.00E-04	2.00E-04	1.00E+00	5.00E+03	5.00E+03	NV	1 00E+00	2.00E+00	1.00E+0
2,4,5-Trichlorophenol	0-0-0	1 00E+01	1 00E-02	2.00E-01	1.00E+00	5 00E+02	5.00E+03	1.00E+03	1 00E-01	2.00E+00	
Dichlorediffuoromethane	0-0-0	1 00E+01	1 00E-02	1.00E+00	4.00E-01	5.00E+01	5.00E+01	NV			1 00E+01
Tetrachlorobonzone	0-0-0	1.00E+04	1 00E-04	2.00E-01	1.00E+00	5.00E+03	5.00E+03		1 00E-01	1 00E+01	4.00E+00
Vinyl Chloride	0-0-0	1 005-04	1 GOE-02	1.00E+00	7.006-04			1.00E+02	1 00E+00	2.00E+03	1.00É+04
				1.002100	7.008-04	5.00E+00	5.00E+00	NV	1.00E+02	1 00E+04	7 00E+00

1,1,1,2   stanchisocothase				11	IRS Pathway Scores		
	Shemical		enioemic	SW/OF/Eav	SW/QW/DWT	SW/OW/HFC	SW/GW/Eav
1,1 - testrachloroschanae		CV2 Mo		•		(Tox/Mob/Per/Bio)	(l:tox/Mob/Pcr/Bio
1,1,2-ternelworethane						<u> </u>	2 00E+00
1,1-inchlorochlane	1,1,2-tetrachloroethane	l *	1	1	1	2 00E-01	2 00h:-0
1,1,2-irichlorochhane	•-•	1				2 00E-01	2 00E+0
1,2-inchlore-thase	,1,2,2-tetrachloroethane		_				2 00E+0
1,2-inchloro-1,2,2-influorochane   75-31-3   2 006-01   NV   NV   NV   NV   NV   NV   NV   N	, i , 2 - trichlorecthans	1		J			N'
1-de: Allorocchane	,1,2-trichloro-1,2,2-triffuoruethane	L L	<del></del>	• • • •			N'
1-deckhorochylene	, I - dichlorocthane	1					N
2,3-inchloropropanae	, I - dichloroethylene		• • •		- 1	1	2 00E-6
2.4 first historopheasean   93-94-3   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-04   2006-0	,2,3-inchiorepropens				•		5 00E+0
2,4-inchlorebeatene	2,4,5-tetrachlorobenzene	95-94-3	:		•		2.00E+0
2dehromon-3-chloropropane	,2,4-trichlorebenzone	120-02-1				•	2:00E*C
2- debrumo-3-chlorogropane	,2-dibromoethans	106-93-4	-			5,552 = 0	., N
2-dechlorocethana   107-08-2   2.008-02   2.008-03   4.008-02   2.008-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04   2.068-04	=	96-12-8	5.00E+05				2 00E+
107-06-2   2.00E-02   2.00E-02   2.00E-02   2.00E-02   2.00E-02   2.00E-02   2.00E-02   2.00E-02   2.00E-02   2.00E-03   NV   NV   NV   NV   NV   NV   NV   N	2-dichlerobeazeas	95-50-1	2.00B+02-			- · · · •	<del>-</del> - ·
2-dichlorocthyleane   78-87-5   2 00E+04   2 00E+02   4.00E+02   2.00E+04   2 00E+03   NV   NV   2.00E+03   3dichlorobrazeae   541-73-1   NV   2 00E+03   NV   NV   2.00E+03   3dichlorobrazeae   541-73-1   NV   2 00E+03   4 00E+03   2.00E+04   2.00E+04   2.00E+04   3dichlorobrazeae   106-44-7   2 00E+04   2 00E+03   4 00E+02   2.00E+04   2.00E+04   3dichlorobrazeae   184-05   106-04   2 00E+04   2 00E+04   4 00E+02   3.00E+04   2 00E+04   3dichlorobrazeae   3.4.6-teztachioropheaol   38 NO 2   5 00E+04   2 00E+04   4 00E+02   3.00E+04   2 00E+04   4.00E+02   3.00E+04		107-06-2	2.00B+02	2.00E+00		•	
78-67-5   2006-02   4.006-02   2.006-03   2.006-04   3.3-dichlorobrazeae   541.73-1   NV   2.006-03   NV   NV   2.006-03   3.3-dichlorobrazeae   541.73-1   NV   2.006-03   4.006-03   2.006-04   2.006-04   3.3-dichlorobrazeae   100.64-7   2.006-02   2.006-03   4.006-02   2.006-00   2.006-04   3.3-dichlorobrazeae   100.64-7   2.006-04   5.006-03   4.006-02   2.006-00   2.006-04   3.3-dichlorobrazeae   100.64-7   2.006-04   5.006-03   1.006-02   5.006-00   5.006-04   4.006-01   2.006-02   2.006-02   4.3-Tr (Salves)   93.72   2.006-03   5.006-03   5.006-04   4.006-01   5.006-03   5.006-03   4.3-Tr (Salves)   93.72   5.006-03   5.006-03   5.006-04   1.006-01   5.006-03   5.006-04   4.3-Tr (Salves)   4.3-Tr (Salves)   4.3-Tr (Salves)   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-03   5.006-	2-dichloroethylens	156-60-5	NA	· 1	- 1	- 1	N
3-dichlorobeazzac		78-87-5	2 00E+04	2 00E+02	*		2 001:+
3-dichloropropeac	•	541 - 73 - 1	NV	2 00E+03	MA	• • • •	2.00E+
4-dichlorobeazene 3.4 o 2 5 00E+02		542 75-6	2 00E:+04	2 00E+03	4 00E+03	• • • • • • • • • • • • • • • • • • • •	2 008+
3,4,6-tetrachlorophenoi    38 %0 2   3.00E+04   3.00E+04   3.00E+02   3.00E		106 46 7	2 00E+02	2 005+03	4 OUE-02	2.00E+00	2.00E+
4.3 TP (Silver) 4.3 TP (Silver) 5.0 TP (Silver		58 90 2	5 00E+04	5 00E+06	1 00E -02	5.00E+00	5.00E+
4.5 trs. blorophenol 4.6 trs. blorophenol 4.6 trs. blorophenol 4.7 trs. blorophenol 4.6 trs. blorophenol 4.7 trs. blorophenol 4.7 trs. blorophenol 4.8 trs. blorophenol 4.9 trs. blorophenol 4.0 trs. blorophenol 4.1 trs. blorophenol 4.2 trs. blorophenol 4.4 da blorophenol 4.5 trs. blorophenol 4.6 da	• • • •	93 72 1	2 006+04	2 00E+06	4 00E-01	2 00E+02	2 00E+
4,0 trx.blurephacol   88-06-2   5.006+03   5.006+03   1.006+01   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03   5.006+03	• • •	95-95-4	5 00E+03	5 00E+06	1 00E-01	5 00E+01	5 00E+
4 D   94-75-7   5 00E+03   5 00E+03   1 00E+00   5 00E+01   5 00E+01   5 00E+01   5 00E+03   5 00E+03   5 00E+04   1 00E+01   5 00E+03   5 00	• •	\$8-06-2	5.006+03	5 00B+07	1.00E+01	5 00E+03	5.00E+
120-83-2   5.00E+05   5.00E+04   1.00E+01   5.00E+03   5.00E+05   1.00E+061   1.00E+061   5.00E+061   1.00E+061	•	94-75-7	5 00E+03	5 00E+03	1 00E+00	5.00E+01	5.00E+
106-89-8   2 00E+03   2 00E+04		120-83-2	5.00E+05	5.00E+04	1 00E+01	5 00E+03	5 00E+
Part	• • • • • • • • • • • • • • • • • • • •		2 00E+03	2.00E+00	4.00E+03	2.00E+03	2.00E+
2-chlorophenol 95-57-8 2 00E+04 2.00E+04 4 00E-01 2.00E+02 5.00E+00 3.3-dichlorophenol 91-94-1 5.00E+04 NV 1.00E-02 5.00E+00 5.00E+00 1.00E-02 5.00E+00 4.00E-02 2.00E+01 NV					1.00E-03	5.00E-01	•
33-36-ichlorobanzidiae   31-94-1   5.00E+04   NV   1.00E-02   5.00E+00   2.00E+00   2.	•	1		2.00E+04	4 00E-01	2.00E+02	2.00€+
S3-32-9   2.00E+03   2.00E+06   4.00E-02   2.00E+01   2.00E+04   NV   NV   NV   NV   NV   NV   NV   N	•		•	U/1100 11		5.00E+00	ŀ
208-96-8	•	1		• • • •		2.00E+01	2.00E+
75-07-0   2.00E+02   2.00E+00   4.00E+02   2.00E+00   3.00E+01   3.00E+02   3.00E+01   3.00E+02				0.000	NV	NV	1
acetons 67-64-1 2.00E+00 2.00E+01 4.00E+00 2.00E+00 2.00E+00 3.00E+01 5.00E+01 5.00E	• •		= :		•	2.00E+02	2 00E+
107-05-8   5.00E+01   5.00E+01   5.00E+02   5.00E+01   5.00E+02   5.00E+01   5.00E+02   5.00E+01   5.00E+02   5.00E+01   5.00E+02   5.00E+01   5.00E+02   5.00E+04	<b>▼</b> .					2.006+00	2 00E+
2.00E+00						•	5.00E-
107-02-8   2.00E+06   2.00E+06   4.00E+03   2.00E+06   2.00E+02   2.00E+03   2.00E+02   2.00E+03   2.00E+02   2.00E+02   2.00E+02   2.00E+02   2.00E+02   2.00E+04		1				2 00E+01	2 OUE+
10   10   10   10   10   10   10   10	•					2.00E+06	2 006
100E+04   100E		1	:		l l		2.00E+
107-13-1   2.00E+04   2.00E+03   4.00E+02   2.00E+04	*					_	1
107-05-1	•	į.		• • • •	• • • • • •		2.00₺•
alpha-BHC         319-84-6         NV         NV         NV         NV           ammionia         7664-41-7         3 50E-02         3.50E-03         7 00E-02         3 50E-02         3 5           antinoc         62-53-3         5 00E+04         5.00E+06         1 00E+04         5.00E+04         5.0	· ·					<del>-</del>	
amenomia 7664-41-7 3 50E-02 3.50E-03 7 00E-02 3 50E-02 3 5 aniline 62-53-3 5 00E+04 5.00E+06 1 00E+04 5.00E+04 5.00E+04	•		=			-	
Americans 62-53-3 5.00E+04 5.00E+06 1.00E+04 5.00E+04 5.00E+04	•			1	1	• • • •	3 50E-
							5 00E+
anth/accec		62-53-3 120-12-7	2 00E+04	3.00E+06 2.00E+07	4.00E-04	2.00E+00	2 006

Chemical				HRS Pathway Scores		
•	CAS No	SW/OF/HPC	SW/OP/Eav	SWIOWIDWT	SW/OW/HFC	SW/OW/I:a
		(Tor/Pers/Bio)	(litux/Pcs/Bio)	(Tox/Mob/Pcs)	(Tow/Mob/Per/Bio)	(laux/Mob/Per/Bio
aroclor 1016	12674-11-2	NV	NV	NV	NV	N'
aroclor 1221	11104-20-2	NV	NV	NV	NV	N'
eroclor 1232	18141-16-5	NV	NV	NV	NV	N'
reclar 1248	12672-29-6	NV	NV	NV	NV	N'
nnecler 1254	11097-49-1	NV	NV	NV	NV	N/
srecler 1260	11006-82-5	NV	NV	NV	NV	,,, ,,,
bonz(a)anthracene	<b>56</b> -55-3	5 00E+07	5 00E+08	1.00E-01	5.00E+03	\$ 00E+0-
CATORS	71-43-2	2 008+05	2 00E+05	4.00E+01	2 00E+05	2 00E+0
cazo(a)pyrcas	50-32-8	5 00E+04	5.00E+08	1.00E+00	5.00E+04	
ionxo(k)fluoranthene	207-08-9	NV	NV	NV	NV	5.00E+0
ouzyl chlorida	100-44-7	3.50E+02	3.50E+02	7.006-02	3.50E+00	/И,
ola-BHC	319-85-7	NV	NV	NV NV		3.50E+00
is (2-chloroethoxyl) methane	111-91-7	5 00E+01	NV	1 00E+02	NV	N/
is (2-ethylhexyl) phthalato	117-01-7	5 00E+04	5.00E+07	1.006-02	5 00E+01	NV
romodichloromethese	75-27-4	2.00E+03	NV		\$ 00E+00	5.00E+03
remoform	75-25-2	NV	NV	4 00E-01	2.00E+01	'NV
romomethane	74 83 9	2 005+03	2 005+04	NV.	ИV	NV
arbon draulfide	75 15 0	2 00E+04		4 00E+00	2 00E+01	2 00E+02
arbon tetrachionde	36 23 3	2 005+04	2 00E+03	4.00E+00	2 00E+02	2 00E+01
Moni	75 87 6	5 00E+03	2 00E+03	4.00E+00	2 00E+02	2.00E+01
hlordage	57-74-9		NV	1 00E+03	5 00E+03	NV
Morobeazeae	108-90-7	\$ 001:+08	5 00E+08	1 00E+00	5 00E+04	5 00E+04
Moreches	75-00-3	3 50E+00	3 50E+01	7 00E-04	3 50E-02	3.50E-01
Moreform	67-66-3	3 50E-03	ИV	7.00E-04	3.50E-03	NV
Moramethane	74-87-3	2 00E+02	2 00E+01	4.00E+01	2.00E+02	2.00E+01
Moromethyl methyl other	''' '''	3 50E-02	3.50E-03	7.00E-03	3 506-02	3.50E-03
resol (mi-cresol) •	107-30-2	3.50E-01	NV	7.00E-01	3.50E-01	NV
rysome	108-39-4	5.00E+01	5.00E+02	1.00E-01	5.006-01	5 00E+00
WA COA	218-01-9	NV	5 00E+06	NV	NV	5 00E+02
yanagas chlorida	98-82-8	2 00E+05	2.00E+04	4.00E+00	2.00E+03	2.00E+02
velohexanone	506-77-4	NV	NV	NV	NV	2.008402 NV
ODD	108-94-1	5.00E+00	5.00E+00	1.00E+00	5.00E+00	5 00E+00
DT	72-54-8	5.00E+06	5.00E+04	1.00E-02	5.00E+02	5 00E+04
chlorodifluoromethene	50-29-3	5.00B+07	5.00E+08	1.006-01	5.00E+03	
ichlorosthyl ether	75-71-8	2.00E+02	NV	4.00E-02	2.00E+00	5.00E+04
chleroisopropyl other	111-44-4	NV	NV	NV	NV	NV
	108-60-1	NV	NV	NV	NV	NV
chloromethyl other eldrin	542-88-1	ИV	NV	NV	NV	NV
ekana phenyl sitrosamine	60-57-1	5 00E+08	5.00E+08	1.006+00	5 00E+04	NV
	86-30-6	5.00E+03	5.00E+04	1.008-01		5.00E+04
schlorohydran hyl costate	106-89-8	NV	NV	NV	5.00E+01	5.00E+02
hyl acctate hyl benzene	141-78-6	2 00E-01	2.00E-01	4 006-01	NV	,NV
, · · · · · · · · · · · · · · · · · · ·	100-41-4	2.00E+02	2 00E+01	4.006-02	2.00E-01	2 00E-01
hylene dibromide	106-93-4	NV	NV	4.00B-02	2.00E+00	2 00E+01
bylene glycol	107-21-1	5 00E-01	5 00E+00	1.00E+00	5 00E-01	· NV

	1		l'	IRS Pathway Scores		
homical	CAS No	SW/OF/HFC	SW/OF/bay	SW/QW/DWT	SW/QW/HFC	SW/GW/Eav
	CV2 MO	(Tox/Pers/Bio)	(biox/Per/Bio)	(Tox/Mob/Pcr)	(Tox/Mob/Pcr/Bio)	(Exox/Mob/Per/Bio
	110-90-5	5 00E+00	5 00E-01	1.00L+01	5 00E+00	5 00E-0
thylene glycol monomethyl ether	1	2 00h+00	2 00E-01	4 001:+00	2 00E+00	2 00E-0
thyl ether	60-29-7	5.006+05	\$ 00E-07	1 00E-02	5 00E+01	5 00E+0
ueranthene	204-44 0	5 00E+05	5 00E+06	1 00E+00	5 001:+03	5 00E+0
luerene	86-73-7	5 00E+00	5 00li-01	1 001:+01	5 00E+00	5 00E+0
ormaldchyde	50-00-0	- 1	5 00E+08	2 00 6+01	1 00E:+05	1 00E+0
epiachlor	76-44-8	\$ 00E+06	5 00E+08	1 00E+04	5 00E+04	5 00E+0
eptachlor epoxide	1024-57-3	5.00E+04		1.00E-01	5 008+02	5 00E+0
exachlorobonzene	118-74-1	5 00E+06	5.006+05	1.00E-01	5 00E+00	5 00li+0
exachlorobutadiene (hexachloro-1,3-butadiene)	67-48-3	5.00E+04	5.00h+07	1 00E+02	\$ 00E+05	5 00b+0
exachlorocyclopontadaene	77-47-4	5.00E+07	5 00E+05	• [	2 00E+03	2 00E+0
cxachloroethane	67-72-1	2.00E+05	2.00E+05	4 00E+00	2 00E+03 5 00E+05	5 00E+C
ne nachloropentadione	77-47-4	5.00E+07	5 00E+05	1 00E+02	2 00E+02	2 00E (
ydrochloric acid	7647-01-0	2.00E+02	2.006-01	4 00E+02	<b>V</b>	2 00E (
nydrogen sulfido	7783-06-4	2.00E+03	2.00E+02	4.00E+03	2 00E+03	
sobutanol	78-83-1	2 00E+00	2 00E+00	4 001:+00	2 00E+00	2 006+
indeno	58-89-9	5 00E+06	5 00E+06	1 00E+01	5 00E+03	\$ 00E+
nethanol	67 56 1	5 006-01	5 00E-01	1 00E+00	5 00E-01	5 00E-0
nethoxychior	72 43 5	5 00E+06	5 00b+08	1.00E-02	5.00E+02	5 00E+
netayl chlorocarbonate	79 22 1	3 00F+01	ил	NV	NV	N
methylene bio(2-chlurusmiline) 4 4	101 14 4	2 00E+05	NV	4 00E 02	2 00E+01	N
pethylae chloride	75 09 2	2 00E+01	2 00E+00	4 00E+00	2.00E+01	2 00b+
metoyi ethyi ke-ans	78 93 3	2 00E+00	2.00E-01	4.00£+00	2 00E+00	2 00₺⊣
met tyl sembutyt ketome	108-10-1	2.005+01	2 00E+00	4 00E+00	2.00E+01	2 00E+
metryl methacerians	80-62-6	2.00E+01	2.00E+00	4.00E+00	2.00E+01	2 00E+
a based	71-36-3	5.00E+01	5.00E+00	1.00E-01	5.00E-01	. \$.00E⊣
asolubulene	91-20-3	2.00E+04	2.00E+05	4 00E-01	2 00E+02	2.00E+
Bijlopevsoue Billopevsoue	98-95-3	5.00E+03	2.50E+01	1.00E+01	5.00E+01	2.50E-
peninchierobenzene	608-93-5	5.00E+06	5.00E+05	1.00E+01	5.00E+04	5.00E+
pentachioroethene	76-01-7	2.00E+02	2.00E+03	4.00E-02	2.00E+00	2.00E+
pentachioroniumbonzens	82-68-8	5.00E+05	NV	1.00B+01	5.00E+03	N
pontacklerophenol	87-86-5	5.00E+04	5.00E+04	1.00E+00	5.00E+02	5.00E+
•	106-47-8	5.00E+03	5.00E+04	1.00E+03	5.00E+03	5.00E+
p-Chloroenilies		NV	2.00E+06	NV	NV	2.00E+
phonothrons	108-95-2	5.00E+00	5.00E+04	1.00E+00	5.00E+00	5.00E+
phonol	129-00-0	5.00E+03	NV	1.00E-02	5.00E-01	N
pyrae itin	110-86-1	5.00E+02	5.00E+01	1 00E+03	5.00E+02	\$.00E+
pyridiae	100-42-5	2.00E+02	2 00E+03	4.00E-02	2.005+00	2.00E+
atyrens	7664-93-9	2.00E+02	2.00E+00	4.005+02	2 008+02	2 00E+
milaric acid	108-88-3	2.00E+02	2.00E+03	4.00E-02	2.00E+00	2 00E+
tolucae	8001-35-2	\$ 00E+07	\$ 00E+08	1 006+01	5.00E+05	\$ 00E+
tonapheas	127-18-4	2.00E+03	2 00E+03	4.00E-01	2 006+01	2 00E+
tetnichloroethene	109-99-9	5 00E+02	NV	1.005-01	5.00E+02	1 002.
tetrahydrofuras	156-60-5	2 00E+03	2 001:+01	4.00E+01	2.00E+03	2 00E+
trans-1,2-dichloroethylene trichloroethylene	79-01-6	2 00E+02	2.00E+03	4.00E-02	2 00E+00	2 00E+

lizhibit I HRS Hazard Data and Pathway Scores for Halogenated Organics

Chancal				HRS Pathway Score		
	CAS No	SW/OP/HPC	SW/OF/Lav	TWD/WD/WS	SW/QW/IIFC	SW/OW/I
		(Tor/Perr/Hun)	(France) Real Real	Carplat /Ban		
trichlating out theory of have		(100010100)	(010112117011)	(10%/Mob/Pct)	(Ton/Mob/Pcr/Bio)	(lisus/Mob/Per/Hou)
The control control control	5-89-4	×	N .	77	AN	
Viayi chlande	73-01-4	3 50E+02	2	7.001-01	3 501:+00	
Aykas (m-xylene) **	104-31-3	2 00E+02	2 00E+04	4 001:-01	2 (M)m	•
Invicac (p-xylcac) **		3 200			* 000000	2 00E+02
24-7-1	200	700000	60E+03	4 00E-02	2 00E+00	10-406-1
a, a - artimetany a puncino i	9	\$ 00E-04	5 00E+04	1 00F±-00	\$ 00£+02	• A
BCatte(s)entanticont	9	5.00E+07	5 00E+04	1 00 H-03	\$ 00E+07	4 Ontine
ways wearys patholese	9	5.00E+03	5.00E+04	005-03	\$ m=:-n=	
Dibear(a,b) anthrocano	9	3.00E+0€	٧٧	1 00 5-00	( ODE O	5
Dimothyl phthelate	7	5.00R+01	\$ 00F-03	1 000 00	, oor o	
Pagrise	}	3			10-300 c	5.00
Matrix ashudada		Onether 7	2	4.00E+00	2.00E+00	
	9	3.50E-03	3.50E-04	74	×-	
	j	5.00H-04	5.00E+05	1.00E+02	· MEL	•
Phihalic sahydride	9	2.00E-01	Z <	A MILLS	3 65 65	9.00
Telume disocyanate	- -	2 MEAN	3 8 6 7	1	10-200.7	
2.3.7.8-Tetrachlerodibenzoinkioria		201100	10-000.7	2	NV.	
2 & Tricklomband	- F	3 00b+07	Z <	1.00E+00	5 00E+03	X X
The second of th		5.00E+03	5.00E+06	1.00E-01	5.00%-01	ŝ
Chemical and Carlo Carlo	7	2 00E+02	× .	A MF-A3	385	J. 94
Tetrachierobcazcao	- - - - -	300		4.000.04	Z.DUE-DUI	
Visyl Chloride		3 00007	CO-ED C	1.00E-00	5 00E+03	5.00E+01
		10-40K K	2			

- The SCDM values for m-cresol were used because no values were available for mixed cresols, m-cresol is more persistant than other isomers.
- •• The SCDM values for sa-xylene and p-xylene were used because no values were available for mixed xylenes. These isomers were used because m-xylene is more bioaccumulative, and p-xylene is more toxic than the other isomers.
- "Tonicity" refers to the HRS human tonicity finess seems [max score = 115-4]
- "OW Mob" refers to the HRS ground-water expliftly factor score based on water solubility and distribution coefficient (Kd) [max score = 1]
- "Air Mob" refers to the air gas mobility factor soom based on the vapor pressure of the gaseous hazardous substitute [max score = 1]
- "Persist" refers to the HRS surface water persistance factor score based on a substance's half-life (based on biodegredation, hydrolysis,
- "PCBroace" refers to the HRS bioaccumulation potential factor score evaluated in the human food chain threat of the surface water pathway ghotolysis, and volatilization) and Kow (stax scere = 1]
- "EnvBioace" refers to the HBS bioaccumulation potential factor score evaluated in the environmental threat of the surface water pathway
- "Ecotox" refers to the HRS ecosystem texticity factor ecors evaluated in the environmental threat of the surface water pathway [mex acory=56+4]
- "GW (Tox/Mob)" refers to the HRS Texicity/Mobility factor score evaluated in the ground-water pathway, and is obtained by multiplying its [max score is life-4]
- "Air (Tou/Mob)" refers to the HRS Toxicity/Mobility factor score evaluated in the air pathway, and is obtained by multiplying its factor factor value componenta (mas acora = 15:4) valus components (max score = 15:4)
- \*SW/OP/DWT (Ton/Pen)\* refers to the Toxicity/Permistence factor ocore evaluated in the overland/flood migration component of the surface water pathway drinking water threat, and is obtained by multiplying its factor value components (max score = 10,000)
- \*SW/OP/HPC (Tox/Pen/Bio)\* refers to the Toxicity/Persistence/Bionecumulation factor neore evaluated in the overland/flood myration max acore=SE+8 component of the surface water pathway human food chain threat, and is obtained by multiplying in factor value components
- "SW/OP/Eav (Bion/Pen/Bio)" refers to the Ecosystem Tonicity/Persistence/Bionecumulation factor value evaluated in the overland/food migration component of the surface water pathway cavironmental threat, and is obtained by multiplying its factor value components max score=SE+8]
- "SW/OW/DWT (Tox/Moh/Put)" refers to the Toxicity/Mobility/Pernistance factor score of the ground water to surface water migration component of the surface water pathway drinking waster threat, and is obtained by multiplying its factor components max scors = 15+4]
- \*SW/OW/HPC (Tou/Mob/Per/Bio)\* refers to the Touicity/Mobility/Persistence/Bioaccumulation factor score evaluated in the ground its factor value components (mex score = 56+6) water to eachee water migratice compensat of the surface water pathways himna food chain threat, and is obtained by multiplying
- "SW/OW/Env (Her/Mob/Par/Die)" raters to the Ecosystem Toxicity/Mobility/Persistence/Bioaccumulation factor ocore evaluated in the ground water to surface water majoration component of the surface water pathway covironmental threat, and is obtained by
- "NY" = so value; "NA" = and applicable

multiplying its factor value components (max score = 56.8)

								Scores or Ka	n <u>k</u> s			****		
Themsest	<del> </del>							HRS Pactor	Scores				<del></del>	
	CAS No	Touchy	GW MOP	Permit	PCBioacc	EnvBloacc	£:cotox	ow	SW/OP/DWT	SW/OP/HI'C	SW/Ol7/Linv	SW/QW/DWT	sw/ow/HPC	SW/OW/E
Lam Lacues	7429-90-5					<del></del>		(Ton/Mob)	(Tox/Pcr)	(Tox/Pers/Bio)	(lidux/Per/Bio)	(Tox/Mob/Pcr)	(Tux/Mub/Per/Hiu)	(Eloz/Mob/Par/t
	1	NV	NV	1	5.00E+01	5 00E+02	1 001:+01	NA	NA	NA	5 00E+03	NA	NA	
rimony.	7440-36-0	1 00E+04	1 00b-02	1	\$.00R-01	2 00F-01	NV	1.00E+02	1 00E+64	5 00E+03	NA	1 001:+02	5 00h+01	
TCRIC	7440-38-2	1 00E+04	1.00B-02	1	\$.00E+00	5.00fi+01	1.00E+01	1.00B+02	1.00E+04	5 00E+04	5 00E+02	1 001:+02	5 00E+02	5 00L
ITUM.	7440-39-3	1.00E+01	1.00B-02	1	5.00H-01	5.00E-01	1.00E+00	1 005-01	1 00E+01	5 00E+00	5 00E 01	1 001: 01	5 00E-02	2
cyllium	7449-41-7	1.00E+04	1.00B-02	1	5.00E+01	5 00E+01	NV	1.00E+02	1 005+04	5 00E+05	NA	1 00E+02	5 00h+03	5.00E
Giorni Lucian	7449-43-9	1.00E+04	2.00B-01	1	5.00E+03	5.00E+03	1 00E+03	2.00E+03	1.00E+04	5 00E+07	5 001:+06	2 00E+03	1 005-07	
	16055-83-1	1.006+00	2.00E-05	0.4	5.00B+04	5 00E+04	1.00E+01	2.00H-05	4.00E-01	2 00E+04	2 001:+05	8.00E-06	1	1 00h
romium	1 1	1.00E+04	NV	1	5.00E+00	5 00E+00	1.008+02	NA	1.00E+04	5.00E+04	5 001:+02	NA NA	4 906-01	4 00E
pper	7440-50-8	NV	1.00B-02	1 ]	5.00B+04	5.00E+04	1.00E+02	NA	NA	NA	5 00E+06	· NA	NA	
ul	7439-92-1	1.00E+04	2.00B-05	1	5.00E+01	5.00B+03	1 00E+03	2.00E-01	1.00E+04	5.00E+05	5 00E+06		NA	5 00E
creury	7439-97-6	1.00E+04	2.00B-05	1	5.00E+04	5.00E+04	1.00E+04	2.005-01	1.00E+04	5 00E+08	5 00E+08	2 00E-01	1 00E+01	1 00h
:kel	7440-02-0	1.00E+04	2.00E-05	1	5.00E-01	5.00E+02	1.00B+01	2.00E-01	1.00E+04	5.00E+03	5 00E+03	2 00E-01	1 001:+04	1 00h
lcaium	7782-49-2	1.00E+02	1.00E-02	1	5.00E+03	5.00E+03	1.00B+02	1.00E+00	1.00E+02	5.00E+05		2 00E-01	1 00E-01	1 00h
ANE	7440-22-4	1.00E+02	2.00E-07	1	5 00E+01	5 00E+01	1.00E+04	2.00E-05	1.006+02		5 00E+05	1.00E+00	5 00E+03	5 00L
llium	7440-28-0	1.00E+03	1.00E-04	1	5.00E+02	5 00E+02	NV	1.006-01	1.00E+03	5 00E+03	5 00E+05	2 00E-05	1 00E 03	1 00h
radium	7440-62-2	1.00E+02	NV	- 1	5 00h 01	3 00b-01	NV	NA	1.00E+02	5.00E+05	NA	1.005-01	5 00E+01	1
c	7440-66-6	1.00E+01	2 00b-03	- 1	3 006-03	5 006-02	1 005+01	2 006-02		5 00E+01	NA	NA	NA NA	1
			· · · · · · · · · · · · · · · · · · ·			1		1 WB-01	1 006+01	5 00E+03	5 00E:+03	2 006-02	1 001:401	1 0012

<sup>&</sup>quot;To acity" refers to the HRS human toxicity factor score [mex ecore = 10,000]

<sup>&</sup>quot;(1% blish" refe a to the HRS ground-water mobility factor score based on water solubility and distribution coefficient (Kd) [max score = 1]

<sup>&</sup>quot;Paraset" refers to the HRS uniface water persentance factor acore based on a substance's half-life (based on biodegredation, hydrolysis, photolysis, and volatization) and Kow [max score = 1]

<sup>\*</sup>PC sees.c \* refers to the HPS beaccumulation potential factor score evaluated in the human food chain threat of the surface water pathway [max score=5E+4]

<sup>&</sup>quot;has Broace" refus to the HRS broaccumulation potential factor score evaluated in the environmental threat of the surface water pathway (max score=5E+4)

<sup>&</sup>quot;ficetox" refers to the HRS ecosystem texicity factor score evaluated in the environmental threat of the surface water pathway [max score = 10,000]

<sup>&</sup>quot;OW (Tox/Mob)" refers to the HRS Texicity/Mebility factor score evaluated in the ground-water pathway, and is obtained by multiplying its factor value components [max score = 10,000]

<sup>&</sup>quot;SW/OP/DWT (Toz/Per)" refers to the Texicity/Persistence factor score evaluated in the overland/flood migration component of the surface water pathway drinking water threat, and is obtained by multiplying its factor value

<sup>&</sup>quot;SW/OF/HPC (Fex/Pers/Bio)" refers to the Texicity/Persistence/Bioaccumulation factor score evaluated in the overland/flood migration component of the surface water pathway human food chain threat, and is obtained by multiplying

<sup>&</sup>quot;SW/OF/Env (Etex/Pers/Bio)" refers to the Ecosystem Toxicity/Persistence/Bioaccumulation factor value evaluated in the overland/flood migration component of the surface water pathway cavironmental threat, and is obtained by

<sup>&</sup>quot;SW/GW/DWT (Tex/Mob/Per)" refers to the Texicity/Mobility/Persistence factor score of the ground water to surface water migration component of the surface water pathway drinking water threat, and is obtained by multiplying its

<sup>&</sup>quot;SW/GW/HPC (fox/Mob/Per/Bio)" refers to the Toxicity/Mobility/Persistence/Bioaccumulation factor score evaluated in the ground water to surface water migration compunent of the surface water pathway buman food chasa threat,

<sup>&</sup>quot;SW/GW/Env (Biox/Mob/Pes/Bio)" refers to the Ecosystem. Toxicity/ Mobility/ Persistance/Bioaccumulation factor acore evaluated in the ground water to surface water migration component of the surface water pathway cavironment "NV" = no value, "NA" = nut applicable

# APPENDIX 9 STATES AND REGIONS IN WHICH TOP 100 RANKED WASTESTREAM COMBINATIONS ARE GENERATED

Rank	RCRA Code	SIC Code	Source Code	Form Code	Volume (tons)	Hazard Score	State	EPA Region
1	D001 F001 F002 F003 F005 U001 U002 U003 U019 U028	2869	A33	B219	14,217	7.09e+13	ΊX	6
2	D002 D006	2833	A32	B207	3,724	3.72e+13	CT	1
3	K022	2869	A33	B606	23,281	2.32c+13	PA	3
4	D001 D002 D019 D032 D033 D034 D039 F002	2869	A33	B219	3,866	1.93c+13	TX	6
5	D001 D007 D008 D018 D022 D026 D027 D028 D033 D036	9999	A99	13219	7,001	1.05c+13	IN	5
6	D001 D002 D003 P002 P020 P024 K017 K018 K020 K028	2821	A37	B219	48,039	4.79c+12	TX	6
7	P003 F005	Unkwn	Unkwn	B219	17,218	2.58c + 12	СТ	1
						i	MA NJ	1 2
							NY	2
	•		, -				PA	3
							VA WV	3
8	D001 D008 F003 F005	4953	A73	B203	4,880	2.43c+12	KY	4
9	D001 D002	2869	A31	B207	3,873	1.93c+12	VA	3
10	Unknown	Unkwn	Unkwn	B206	84,191	1.68c+12	СТ	1
							RI	1
							DE DE	3
						٠.	MD	3
		1 -				·	PA	3
	•						VA IL	3
							ΤX	6
					ŀ		IA	7
							KS MO	7 7
							CA	9
11	D001 D002	2869	A35	B219	95,042	1.42c+12	17X	6
12	D001 D018 D019 D039 F024	2869	A74	B202	26,708	1.33c+12	I.A	- 6
13	D001 D028 F037 F038	2911	A89	B205	6,785	6.77c+11	TX	6
14	K002	2865	A33	B203	6,554	6.54c+11	LA	6
15	D001 D005 D006 D007 D008 D018 D026 D035 F001 F002	Unkwn	Unkwn	B219	31,348	6.26e+11	MI	5
16	D019 D022 D032 D039 D043 K018 K020	2869	A33	B219	3,132	6.25c+11	ТX	6
17	D001 D002 U008 U113	2869	A33	B101	189,524	5.67c+11	ΤX	6

Rank	RCRA Code	SIC Code	Source Code	Form Code	Volume (tons)	Huzurd Score	State	EPA Region
18	D001 K013 U003	2869	A33	B219	5,554	5.54c+11	ΊX	6
19	D001 F001 F002 F003	Unkwn	Unkwn	B204	10,782	5.38c+11	FL KY IN OH WI IA	4 4 5 5 5 7
20	D001 D008	2821	A33	B602	13,395	5.35e+11	KY	4
21	K048 K049 K051	Unkwn	Unkwn	Unkwn	3,393	5.08c+11	OH LA	5
22	F003 F005	2819	A	В	6,101	4.57c+11	TN	4
23	12001 F003 F005	Unkwn	Unkwn	B203	5,692	4.54c+11	CT MA ME NJ NY MD PA VA AL FL MS NC SC TN IN	1 1 2 2 3 3 3 4 4 4 4 4 4 5
							MI MN OH I.A KS MO CO ND CA	5 5 5 6 7 7 8 8

Rank	RCRA Cude	SIC Code	Source Code	Form Code	Volume (tons)	Huzard Score	State	EPA Region
24	D001 D004 D005 D006 D007	Unkwn	Unkwn	B204	22,251	4.44c+11	TN IN MI	4 5 5
25	D001 D002 F002 F003 F005 U002 U012 U031 U044 U080	2834	A37	H201	4,163	4.15c+11	GA	4
26	D001 D002 D003 D006 D018 D023 D024 D025 D026	2869	A33	B219	18,825	3.76c+11	TX	6
27	D001 D018 D019 D022 D028	2869	A37	H202	8,416	3.36c+11	LA	6
28	KOSI	2911	A89	B603	6,217	2.92c+11	TX	6
29	D001 D002 D007 D018 D021 P002 F003 F005	2865	A31	B204	4,781	2.86c+11	AR	6
30	D001 D002 D003 D018 D026 D035 F002 F003 F004 F005	2869	A33	B219	14,194	2.83e+11	ТX	6
31	D001 D004 D005 D006 D007 D008 D009 D010 D011 D016	Unkwn	Unkwn	B202	6,591	2.63c+11	МО	7
32	D001 F003 F005	Unkwn	Unkwn	B204	8,747	2.62c+11	NJ VA WV KY IL IN	2 3 3 4 5 5
					-	# 	MI OH MO NE CA	5 5 7 7 9
33	D001 D002	2819	A37	B219	15,997	2.39c+11	TX	6
34	K049	2911	A75	B202	3,316	2.32c+11	II.	5
35	D001 D004 D005 D006 D007 D008 D009 D010 D011 D018	Unkwn	Unkwn	B202	5,565	2.22c+11	AL TN CA	4 4 9
36	D001 D005 D006 D007 D008	7389	A89	B204	10,883	2.17c+11	TX	6
37	D018 F037 F038 K048 K049 K050 K051	2911	A75	B603	10,580	2.11c+11	I.A	6
38	D001 D004 D005 D006 D007 D008 D009 D010 D016 F001	7389	A71	B219	4,743	1.89c+11	TX	6
39	D001 D002 D003 D004 D005 D006 D007 D008 D009 D010	4953	A99	B114	4,564	1.82c+11	TX	6
40	D001 D004 D005 D006 D007 D008 D010 D011 D018 D035	2899	A89	B204	4,531	1.81c+11	AL	4
41	D001 D002 D007 D018 D021 F002 F003 F005	2865	A34	B204	4,316	1.72c+11	AR	6

Rank	RCRA Code	SIC Code	Source Code	Form Code	Volume (tons)	llazurd Score	Stute	EPA
42	D001 D007 D008 D018		<del></del>					Region
43		2911	A89	B204	8,564	1.71c+11	OH	5
4.5	D001 D005 D006 D007 D008	Unkwn	Unkwn	B407	7,826	1.56c+11	MI AR	5
44	D001 D018 K048 K049	2911	Unkwn	B204	3,669	1.46c+11	ОН	5
45	Unknown	3221	A54	B206	7,914	1.18c+11	NJ	2:
46	D001 F024	2819	A33	B219	14,893	1.04e + 11	LA	6
47	K048	2911	A75	B503	19,996	9.98c + 10	TX	6
48	D001 D004 D005 D006 D007	Unkwn	Unkwn	B407	4,509	9.00c+10	GA	<del> </del>
		0			4,505	3.00C+10	TX	6
49	D001 D002	2869	A33	B219	7,412	8.87c+10	VA	3
				•			ΤX	6
50	D005 D006 D008 F001	4953	Unkwn	H204	4,348	8.68c + 10	OH	5
51	D018 D038 K022 K083	2865	Α	В	17,303	8.63c + 10	ОН	5
-52	D001 D005 D006 D007	Unkwn	Unkwn	B204	3,775	7.53c+10	OH	5
53	D001 D011 D018 D021 D022	3861	A49	B204	9,390	7.49c+10	NY	2
54	D001 F001 F003 F005	Unkwn	Unkwn	B204	5,956	7.43e+10	RI	1
							NY	2
			ļ				II. Mi	5
			1				OH	5
55	D001 D002 D007		<u> </u>				NM	6
56		2869	A33	B602	36,709	7.32c+10	TX	6
57	D001 D005 D006 D007 D008 F001 F002 F003 F004 F005	7389	A71	B206	3,518	7.02c + 10	ΤX	6
	D001 D005 D006 D007 D008 F003 F005	2821	A73	B602	3,410	6.80c + 10	WI	5
58	D001 D002 D005 D006	4953	Unkwn	B204	3,295	6.57c+10	OH	5
59	D001 D002 D019 D022 D027 D028 D029 D032 D033 D034	2869	A99	B494	6,435	6.42c+10	TX	6
60	D001 D006 D008 P002	Unkwn	Unkwn	B403	3,168	6.32c+10	ОН	5
61	F001 F002 F003 F005		<del> </del>		<del>                                     </del>		TX	6
		Unkwn	Unkwn	B204	6,975	6.26c+10	PR	2
	-					1	FL IN	4
				1	1		TX	5

Renk	RCRA Code	SIC Code	Source Code	Form Code	Volume (tons)	Hazard Score	State	EPA Region
62	D001 D005 D006 D007 D008 D011 D022 D035 D039 F001	Unkwn	Unkwn	Unkwn	3,124	6.23c + 10	OH	5
63	D001 D002 D007 D008 D018 D035 F001 F003 F005 U009	2869	A37	B219 ·	5,679	5.67e+10	ΤX	6
64	D001 F001 F002 F005	Unkwn	Unkwn	B204	10,929	5.45e+10	WI	5
•					ž.		Q	11
65	D008	Unkwn	Unkwn	Unkwn	5,357	5.34e + 10	MA	1
		-					VT	1
		· ·					NY FL	4
	· · · · · · · · · · · · · · · · · · ·			1			KY	4
							NC	4
		į				,	SC	4
							TN IL	5
		14 /4 4		1			IN	5
			1.	<b>!</b> ;	•	·	MI	5
		1					MN	5
							ОН	5
		,					i.a OK	6
							TX	6
							KS	7
-		e.					MO	7
			1			·	NE	7
				3	•		AZ CA	9
				1			o	l ú.
66	D001 F004	2821	A33	B602	3,990	5.18e+10	NY	2
67	D001 D007 D008 F001 F002 F003 F005	9999	Unkwn	B202	4,866	4.85c+10	TN	4
68	K022	2865	A33	B602	9,432	4.71c+10	TX	6
69	K022	2869	A33	B219	10,846	3.25c+10	TX	6
70	D001 D002 D007	2869	A33	B219	16,099	3.21c+10	TX	6
71	D001 F001 F002 F003	2899	A89	B204	5,825	2.91c+10	AL	- 4
72	D001 D002 F003 F005 K038 P094	2879	A37	B101	35,136	2.80c+10	MO	7
73	D001 D002 D003 D004 D005 D006 D007 D008 D009 D010	2869	A33	B105	13,182	2.63c+10	ΤX	6
74	K017 K019 K020	2869	A33	B601	13,073	2.61c+10	ΤX	6

Rank	RCRA Code	SIC Code	Source Code	Form Code	Volume (toms)	Hazard Score	State	EPA Region
75	D001	2512	A92	B403	4,322	2.59c+10	NC	4
76	D001 D018 D043 F001 F002 F003 F004 F005	Unkwn	Unkwn	H204	15,509	2.48c+10	VA	3
77	F001 F002 F003	Unkwn	Unkwn	Unkwn	4,822	2.41c+10	NJ GA TN IN OH TX CO	2 4 4 5 5 6 8
78	K022	2865	A35	B602	4,609	2.30c+10	1X	6
79	D001 F003 F005	3053	A56	B403	3,465	2.28c+10	VA	3
80	D002 D021 D028 P003 F005	2879	A37	B101	130,948	2.22c+10	MO	7
81	D001	2869	A35	13207	10,732	2.14c+10	TX	6
82	D001	2869	A35	H606	3,175	1.27c+10	MI	5

Ronk	RCRA Code	SIC Code	Source Code	Form Code	Volume (tons)	liuzurd Score	State	EPA Region
83	D001	Unkwn	Unkwn	Unkwn	25,371	1.01c+10	CT MA RI VT NJ NY	1 1 1 2 2,
-							PA VA WV AL FL GA	3 3 4 4 4
		·					KY MS NC SC TN IL	4 4 4 4 5
-						•	IN MI MN OH WI AR	5 5 5 5 5
							LA TX IA KS MO NE	6 7 7 7 7
							MT WY AZ CA NV WA	8 8 9 9 9
							FC Q	11 11

Rank	RCRA Code	SIC Code	Source Code	Form Code	Volume (tons)	Hazard Score	State	EPA Region
			<del></del>	<del></del>	+		<del> </del>	
84	10001 F001 F002 F003 FX05 	Unkwn	Unkwn	B202	5,922	9.45c+09	II.	5
							IN	5
							ОН	5
	Fact Days	<del>- </del>	<b> </b>	-			WI	5
85	D001 D002	Unkwn	Unkwn	Unkwn	3,477	6.94c+09	MD KY	3
			Ī				TN	
			1	,			IL	5
			İ				OH WI	5
							ΪÀ	6
							OK	6
							TX CA	6
86	D001 F002 F003 F005	2384	A37	B101	18,747	5.61c+09	PR	2
<u></u>				10.01	10,747	3.010 +03	IN	5
87	1901 1902 1903 1 005	2833	A35	B101	26,284	2.10e+09	IN	5
MH.	EMD1 13022	2869	A37	B202	3,414	1.16c+09	LA	6
H4	H02 H05	2834	A37	B101	28,640	1.14c+09	IN	5
9()	D001 D002 D003	2879	A37	B102	27,247	8.16c+08	TX	6
91	D002	2869	A09	B207	8,001	7.98c+08	TX	6
92	D001	2869	A35	B219	12,842	6.41c+08	wv	3
			ļ				AL	4
	•	1					I.A TX	6
93	D001 D002 D003 D018 D021 D023 D024 D025 D026 D035	2869	A33	B219	7,418	5.03e+08	17	<del> </del>
94	K027	2865	A33	B409	11,123	4.44c+08	LA	6
95	D001 D002 D003 D005 D018 D021 D023 D024 D025 D026	2869	Unkwn	Unkwn	6,075	4.12c+08	OH	
96	D001 D002 F003	2819	A	В	5,323	1.06c+08	TN	5
97	K027	2865	A33	B403	4,457	8.89c+07	WV	4
98	D001 F002 F003	2833	A35	B101	18,154	5.43c+07		3
99	F002	2879	A35	B101	37,447	1.79c+07	IN CA	9

Kank	RCRA Code	SIC Code	Source Code	Form, Code	Volume (tons)	Hazard Score	State	EPA Region
100	F002 F003 F005	2384	A37	B101	6,414	5.12e+06	PR	2

